

**Falcon Refinery Superfund Site
Ingleside
San Patricio County, Texas
TXD 086 278 058**

Monthly Progress Report # 61

May 2011

Prepared for

**National Oil and Recovery Corporation
3717 Bowne Street
Flushing, NY 11354**

Prepared by



**505 East Huntland Drive
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June 9, 2011

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1.0 INTRODUCTION

This sixty-first Monthly Progress Report is submitted in accordance with the Falcon Refinery Site Administrative Orders on Consent for Removal Action and Remedial Investigation / Feasibility Study between the U.S. Environmental Protection Agency (U.S. EPA) and National Oil Recovery Corporation (NORCO).

This Monthly Progress Report and subsequent reports will address activities associated with both of the orders.

The next monthly progress report, covering June, 2011 will be submitted on or before July 10, 2011.

2.0 COMPLETED ACTIVITIES

2.1 Removal Action Activities

During May the volume of liquid waste was measured in each of the above ground storage tanks (Figure 1) and samples were obtained to characterize the waste. Provided as Table 5 are the results of the tank gauging, which shows that a majority of the liquid waste is currently in Tanks 26 and 30. During June of 2009 measurements indicated that liquid waste had been removed from the tanks, as a result the current liquid is comprised of rainwater that fell into the tanks since June 2009 and came in contact with varying amounts of sludge that remained in the tanks.

Each of the liquid samples from the tanks was analyzed for volatile organic compounds (VOC), semi-volatile compounds (SVOC), metals and pH (Appendix A). Results of the sampling indicated that the rainwater that fell into the tanks is impacted minimally by the sludge that remained in the tanks. As an example, the analytical results from Tank 27, which contains approximately 210,193 gallons of liquid, met all TCEQ Texas Risk Reduction Program (TRRP) Protective Concentration Levels (PCL) for residential drinking water standards.

Provided as Table 6 is a compilation of all the analytical results, from the May sampling, that exceeded the respective PCLs for residential drinking water.

To date a total of approximately 7,774,721 gallons of hazardous waste have been removed from all of the above ground tanks and disposed via deep well injection at Texas Molecular.

Prior to the beginning of liquid waste disposal in October 2004, the volume of waste in the above ground storage tanks was measured at 6,844,094 gallons. Apparently due to holes in the tops of the tanks the volume of waste has increased due to rainfall, since more waste has been disposed of than was originally measured.

A compilation of hazardous liquid waste disposal is included as Table 1.

2.2 Remedial Investigation / Feasibility Study (RI/FS)

NORCO will implement the Phase II RI/FS Field Sampling Plan when approval is granted by the EPA.

3.0 CHANGES MADE IN THE PLANS DURING THE REPORTING PERIOD

Removal Action Work Plan Addendum No. 3 (Appendix B), which deals primarily with the disposal of the remaining waste in the tanks, was submitted to the EPA.

4.0 COMMUNITY RELATIONS

The EPA has developed a web site to display information about the Removal Action and RI/FS activities. Information can be found by going to www.epaossc.net and selecting web sites, then Region 6 and then the Falcon Refinery Site.

5.0 CHANGES IN PERSONNEL DURING THE REPORTING PERIOD

No changes were made during May.

6.0 LIST OF PROJECTED WORK FOR THE NEXT TWO MONTHS

6.1 Removal Action Work projected for the next two months includes:

- Implement Removal Action Work Plan Addendum No.3;
- Pending EPA approval move all liquid into Tank 30;
- As tanks are emptied and cleaned, determine if the tanks will be repaired or removed;
- Continued site maintenance; and
- Apply for a land discharge permit through the TCEQ.

6.2 RI/FS Work projected for the next two months includes:

- Implementing the Phase II Field Sampling Plan.

7.0 LABORATORY / MONITORING DATA

Analytical results from the sampling of the above ground tanks is provided in Appendix B.

FIGURE



N



0 200 400
Feet

ABOVE GROUND STORAGE TANK MAP

FALCON REFINERY
INGELSIDE, SAN PATRICIO COUNTY, TEXAS

PROJECT NO.: 182978

DATE: 4/29/2011



505 EAST HUNTLAND DRIVE
SUITE 250
AUSTIN, TEXAS 78752
512-329-6080

FIGURE
1

Source: National Agriculture Imagery Program
(NAIP) 2009 Aerial Photography.

TABLES

Table 1. Hazardous Liquid Waste Disposal

| DISPOSAL FACILITY | ADDRESS | PHONE NO. | EPA ID NO. | CONTACT |
|---|---|---------------------|--------------|---------------------|
| Texas Molecular Corpus Christi Services, LP | 6901 Greenwood Dr. Corpus Christi, TX | 361-852-8284 | TXR000001016 | Robert Rodriguez |
| RQ, HAZARDOUS WASTE LIQUID N.O.S., 9, UN3082, III (D007, D008, D018) | | | | |
| | Month | Volume (gal) | | |
| | October-04 | 53,832 | | |
| | November-04 | 734,763 | | |
| | December-04 | 879,158 | | |
| | January-05 | 783,881 | | |
| | February-05 | 551,444 | | |
| | March-05 | 565,489 | | |
| | April-05 | 445,107 | | |
| | May-05 | 471,311 | | |
| | December-05 | 42,550 | | |
| | January-06 | 58,740 | | |
| | February-06 | 59,140 | | |
| | March-06 | 0 | | |
| | April-06 | 29,371 | | |
| | May-06 | 59,018 | | |
| | June-06 | 97,151 | | |
| | July-06 | 118,743 | | |
| | August-06 | 148,509 | | |
| | September-06 | 109,908 | | |
| | October-06 | 86,665 | | |
| | November-06 | 140,498 | | |
| | December-06 | 85,813 | | |
| | January-07 | 118,541 | | |
| | February-07 | 107,985 | | |
| | March-07 | 152,493 | | |
| | April-07 | 121,588 | | |
| | May-07 | 150,368 | | |
| | June-07 | 87,900 | | |
| | July-07 | 143,485 | | |
| | August-07 | 94,727 | | |
| | September-07 | 0 | | |
| | October-07 | 50,298 | | |
| | November-07 | 151,227 | | |
| | December-07 | 112,285 | | |
| | January-08 | 119,353 | | |
| | February-08 | 88,777 | | |
| | March-08 | 60,913 | | |
| | April-08 | 18,695 | | |
| | May-08 | 25,349 | | |
| | June-08 | 0 | | |

Table 3. Contaminated Soil and Oily Debris Disposal

| DISPOSAL FACILITY | ADDRESS | PHONE NO. | EPA ID NO. | CONTACT |
|--|------------------------------|--------------------|--------------|-------------------|
| U.S. Ecology Texas L.P. | P.O. Box 307 Robstown, TX | 361-387-3518 | TXD069452340 | Glenda Felkner |
| PETROLEUM CONTAMINATED SOIL AND OILY DEBRIS | | | | |
| | | | | |
| | Month | Volume (cy) | | |
| | October-04 | 0 | | |
| | November-04 | 0 | | |
| | December-04 | 40 | | |
| | January-05 | 0 | | |
| | February-05 | 0 | | |
| | Total | 40 | | |
| RQ, HAZARDOUS WASTE SOLID, N.O.S., LEAD, 9 NA3077, PGIII (OILY SLUDGE AND SOIL) | | | | |
| | | | | |
| | Month | Volume (cy) | | |
| | February-05 | 15 | | |
| | Total | 15 | | |

Table 4. Oil and Filter Disposal

| DISPOSAL FACILITY | ADDRESS | PHONE NO. | EPA ID NO. | CONTACT |
|---|--|---------------------|--------------|---------------------|
| Texas Molecular Corpus Christi Services, LP | 6901 Greenwood Dr Corpus Christi, TX | 361-852-8284 | TXR000001016 | Robert Rodriguez |
| RECYLCED OIL AND FILTERS | | | | |
| | | | | |
| | Month | Volume (gal) | | |
| | January-05 | 403 | | |
| | February-05 | 0 | | |
| | Total | 403 | | |
| DISPOSAL FACILITY | ADDRESS | PHONE NO. | EPA ID NO. | CONTACT |
| Midstate Environmental Services, LLC | 2203 Tower Road Robstown, TX | 361-387-2171 | TXR000051227 | Lloyd Cooke |
| RECYLCED OIL AND FILTERS | | | | |
| | | | | |
| | Month | Volume (gal) | | |
| | January-05 | 16,651 | | |
| | February-05 | 0 | | |
| | Total | 16,651 | | |

Table 5. Falcon Refinery Tank Gauging

| <u>Tank ID</u> | <u>Radius (ft)</u> | <u>Fluid Height (ft)</u> | <u>Tank Height (ft)</u> | <u>Estimated Gallons of Liquid</u> |
|-----------------------|---------------------------|---------------------------------|--------------------------------|---|
| 27 | 53.89 | 3.08 | 40 | 210,193 |
| 26 | 53.89 | 7.41 | 40 | 505,692 |
| 10 | N/A | not measured | 40 | unknown |
| 2 | 25 | 7.25 | 24 | 106,480 |
| 7 | 23.66 | 7.5 | 32 | 98,660 |
| 30 | 85 | 3.5 | N/A | 594,234 |
| 24 | 16.66 | 1 | 32 | 6,522 |
| 20 | 16.66 | 32.16 | 32 | 209,758 |
| 22 | 16.66 | 0 | 32 | 0 |
| 23 | 16.66 | 0.29 | 32 | 1,891 |
| 19 | 16.66 | 0 | 32 | 0 |
| 18 | 16.66 | 0.375 | 32 | 2,446 |
| 21 | 16.66 | 4.5 | 32 | 29,350 |
| Y1 | 9.5 | 6 | 10.25 | 12,725 |
| 17 | 16.66 | 1 | 32 | 6,522 |
| n2 | 6 | 5 | 15.41 | 4,230 |
| n1 | 6 | 15 | 15.41 | 12,690 |
| X1 | 9 | 2 | 21 | 3,807 |
| X2 | 9 | 2 | 21 | 3,807 |
| X3 | 7 | 2 | 16 | 2,303 |
| Total Gallons | | | | 1,801,393.52 |

Table 6. Falcon Refinery May 2011 Analytical Results

| Groundwater Analytical Results Falcon Refinery Superfund Site Ingleside, San Patricio County, Texas | | Sample ID | TANK 26 | TANK 10 | TANK 30 | TANK 7 | TANK 20 | TANK 2 | TANK 27 |
|---|---------------------------------|-------------------------------------|-----------|-----------|-----------|-----------|-----------|-----------|-----------|
| | | Laboratory | TestAmer | TestAmer | TestAmer | TestAmer | AnalSys | AnalSys | AnalSys |
| | | Date Collected | 5/20/2011 | 5/20/2011 | 5/20/2011 | 5/20/2011 | 5/10/2011 | 5/10/2011 | 5/10/2011 |
| Analyte | TRRP Tier 1 PCL ¹ | Residential Source Area, Class 1 | mg/L | mg/L | mg/L | mg/L | mg/L | mg/L | mg/L |
| VOCs (8260) | CAS | GW/GW _{Ing} | | | | | | | |
| 1,3,5-Trimethylbenzene | 108-67-8 | 1.222098214 | 0.012 | 0.0012 | 0.068 | 0.021 | LR | LR | LR |
| 1,2,4-Trimethylbenzene | 95-63-6 | 1.222098214 | 0.044 | 0.0013 | 0.19 | 0.081 | LR | LR | LR |
| 2-Butanone | 78-93-3 | 15 | LR | LR | LR | LR | 0.0854 | LR | LR |
| 4-Methyl-2-pentanone | 108-10-1 | 1.955357143 | LR | LR | LR | LR | 0.00621 | LR | LR |
| Acetone | 67-64-1 | 21.99776786 | LR | 0.015 | LR | LR | 0.624 | LR | LR |
| Benzene | 71-43-2 | 0.00500 | 0.32000 | 0.00730 | 2.50000 | 0.24000 | 0.00778 | 0.19900 | LR |
| Ethylbenzene | 100-41-4 | 0.70000 | 0.13000 | 0.00120 | 0.29000 | 0.17000 | 0.00333 | 0.16600 | LR |
| Methyl tert-butyl ether | 1634-04-4 | 0.24442 | 0.78000 | 0.00170 | LR | 1.50000 | 0.13300 | 0.29000 | 0.00843 |
| Naphthalene | 91-20-3 | 0.48884 | 0.21000 | LR | 0.18000 | 0.37000 | LR | 0.01430 | LR |
| o-Xylene | 95-47-6 | 10.00000 | LR | LR | LR | LR | 0.00159 | 0.03050 | LR |
| Styrene | 100-42-5 | 0.10000 | LR | LR | LR | LR | 0.00229 | LR | LR |
| Toluene | 108-88-3 | 1.00000 | 0.05000 | 0.00070 | LR | 0.51000 | 0.00279 | 0.02630 | LR |
| Xylenes | 1330-20-7 | 10.00000 | 0.25000 | 0.00440 | 1.40000 | 0.53000 | 0.00241 | 0.43200 | LR |
| SVOCs (8270) | CAS | GW/GW _{Ing} | | | | | | | |
| 2,4-Dimethylphenol | 105-67-9 | 0.48884 | LR | LR | 0.11000 | 0.06600 | LR | 0.02060 | LR |
| 1-Methylnaphthalene | 90-12-0 | 0.03147 | LR | LR | LR | LR | LR | 0.01090 | LR |
| 2-Methylnaphthalene | 91-57-6 | 0.09777 | LR | LR | 0.20000 | 0.22000 | LR | LR | LR |
| 2-Methylphenol | 95-48-7 | 1.22210 | LR | LR | LR | LR | 0.13800 | LR | LR |
| 4-Methylphenol | 106-44-5 | 0.12221 | LR | LR | LR | LR | 0.16500 | LR | LR |
| Bis(2-ethylhexyl)phthalate | 117-81-7 | 0.00600 | LR | LR | LR | 0.08900 | LR | LR | LR |
| Chrysene | 218-01-9 | 0.12500 | LR | LR | LR | LR | 0.01720 | LR | LR |
| Phenanthrene | 85-01-8 | 0.73326 | LR | LR | LR | 0.17000 | 0.01470 | LR | LR |
| Phenol | 108-95-2 | 7.33259 | LR | LR | LR | LR | 0.50900 | LR | LR |
| Pyrene | 129-00-0 | 0.73326 | LR | LR | LR | LR | 0.01870 | LR | LR |
| Metals (6010/7470) | CAS | GW/GW _{Ing} | | | | | | | |
| Arsenic | 7440-38-2 | 0.01000 | LR | LR | LR | LR | 0.01190 | 0.00338 | LR |
| Barium | 7440-39-3 | 2.00000 | 0.75000 | 0.18000 | 3.60000 | 0.88000 | 0.00540 | 0.47400 | 0.20900 |
| Cadmium | 7440-43-9 | 0.00500 | LR | LR | LR | LR | 0.00207 | LR | LR |
| Chromium (total) | 7440-47-3 | 0.10000 | LR | LR | LR | 0.01600 | LR | 0.00491 | LR |
| Lead | 7439-92-1 | 0.01500 | 0.02200 | 0.02800 | LR | 0.12000 | 0.05360 | 0.02520 | LR |
| Mercury | 7439-97-6 | 0.00200 | LR | LR | LR | LR | 0.00108 | LR | LR |
| Miscellaneous | | | | | | | | | |
| pH | -- | NL | LR | LR | LR | LR | 13.00000 | LR | LR |
| Notes: | | | | | | | | | |
| ¹ 30 TAC 350.51 | | | | | | | | | |
| ² Texas Secondary Drinking Water Standard | | | | | | | | | |
| NL - Not Limit (no PCL) | | | | | | | | | |
| LR - Less than reporting limit (RL) | | | | | | | | | |
| Data Qualifiers: | | | | | | | | | |
| Concentration exceeding Critical PCL. | | | | | | | | | |

APPENDIX A

Client: Gainco, Inc.
Attn: Paul Supak
Address:

Phone: FAX:

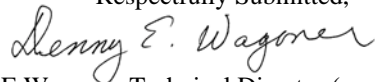
Report#/Lab ID#: 352492 Report Date: 05/16/11
Project ID: Falcon Refinery
Sample Name: Tank 20
Sample Matrix: water
Date Received: 05/10/2011 Time: 16:11
Date Sampled: 05/10/2011 Time: 10:30

REPORT OF ANALYSIS

QUALITY ASSURANCE DATA ¹

| Parameter | Result | Units | RQL ⁵ | Blank | Date | Method ⁶ | Data Qual. ⁷ | Prec. ² | Recov. ³ | CCV ⁴ | LCS ⁴ |
|--|---------|----------|------------------|---------|----------|---------------------|-------------------------|--------------------|---------------------|------------------|------------------|
| A/BN extraction-625/8270 | --- | --- | --- | --- | 05/13/11 | 3510 | --- | --- | --- | --- | --- |
| Metals Dig.-Hg | --- | --- | --- | --- | 05/11/11 | 7470 & SM3112B | --- | --- | --- | --- | --- |
| Metals Dig.-Total | --- | --- | --- | --- | 05/12/11 | 200.2 & 3005A | --- | --- | --- | --- | --- |
| pH (@T=21.9°C) | 13 | pH units | --- | --- | 05/11/11 | 9040C&SM4500HB | --- | 0 | -NA- | 100 | -NA- |
| Arsenic/ICPMS | 0.0119 | mg/L | 0.002 | <0.002 | 05/13/11 | 6020A & 200.8 | --- | 0 | 99.5 | 98.3 | 99.6 |
| Barium/ICPMS | 0.0054 | mg/L | 0.002 | <0.002 | 05/13/11 | 6020A & 200.8 | --- | 1 | 104.9 | 106.6 | 100 |
| Cadmium/ICPMS | 0.00207 | mg/L | 0.001 | <0.001 | 05/13/11 | 6020A & 200.8 | --- | 0.8 | 93.9 | 100.5 | 92.7 |
| Chromium/ICPMS | <0.002 | mg/L | 0.002 | <0.002 | 05/13/11 | 6020A & 200.8 | J, | 1.5 | 99.5 | 98.5 | 100 |
| Lead/ICPMS | 0.0536 | mg/L | 0.002 | <0.002 | 05/13/11 | 6020A & 200.8 | --- | 0.1 | 87 | 95.6 | 92.2 |
| Mercury/CVAA | 0.00108 | mg/L | 0.0002 | <0.0002 | 05/12/11 | 7470&SM3112B | --- | 3.08 | 90 | 107 | 107 |
| Selenium/ICPMS | <0.002 | mg/L | 0.002 | <0.002 | 05/13/11 | 6020A & 200.8 | J, | 1.7 | 106.2 | 104 | 106.9 |
| Silver/ICPMS | <0.001 | mg/L | 0.001 | <0.001 | 05/13/11 | 6020A & 200.8 | --- | 1 | 92.7 | 97.6 | 92.2 |
| Volatile organics-8260 | --- | --- | --- | --- | 05/13/11 | 8260b(5030/5035) | --- | --- | --- | --- | --- |
| 1,1-Dichloroethane | <1 | µg/L | 1 | <1 | 05/13/11 | 8260b & 624 | --- | 2.8 | 138 | 128.1 | 125.9 |
| 1,1-Dichloroethene | <1 | µg/L | 1 | <1 | 05/13/11 | 8260b & 624 | --- | 2.4 | 106.7 | 98.3 | 99.2 |
| 1,1,1-Trichloroethane | <1 | µg/L | 1 | <1 | 05/13/11 | 8260b & 624 | L,S1, | 2.4 | 139.6 | 126.9 | 125.6 |
| 1,1,1,2-Tetrachloroethane | <1 | µg/L | 1 | <1 | 05/13/11 | 8260b & 624 | --- | 2.1 | 105.6 | 101.5 | 103.4 |
| 1,1,2-Trichloroethane | <1 | µg/L | 1 | <1 | 05/13/11 | 8260b & 624 | --- | 2.7 | 121.4 | 119.4 | 120.5 |
| 1,1,2,2-Tetrachloroethane | <1 | µg/L | 1 | <1 | 05/13/11 | 8260b & 624 | C, | 0.5 | 126.9 | 136.6 | 129.9 |
| 1,2-Dibromo-3-chloropropane | <2 | µg/L | 2 | <2 | 05/13/11 | 8260b & 624 | --- | 3.8 | 107 | 105.4 | 106.1 |
| 1,2-Dibromoethane (Ethylene dibromide) | <1 | µg/L | 1 | <1 | 05/13/11 | 8260b & 624 | --- | 3.8 | 115.2 | 114.1 | 112.5 |

This analytical report is respectfully submitted by AnalySys, Inc. The enclosed results reflect only the sample identified above. The results have been carefully reviewed and to the best of my knowledge, unless otherwise indicated, meet NELAC requirements as described by AnalySys, Inc.'s Quality Assurance/Quality Control Program. © Copyright 2003, AnalySys, Inc., Austin, TX. All rights reserved. No part of this publication may be reproduced or transmitted in any form or by any means without the express written consent of AnalySys, Inc.

Respectfully Submitted,

D.E. Wagoner, Technical Director (or designee)

1. Quality assurance data for the sample batch which included this sample. 2. Precision (PREC) is the absolute value of the relative percent difference between duplicate results. 3. Recovery (Recov.) is the percent of analyte recovered from a spiked sample. 4. Calibration Verification (CCV) and Laboratory Control Sample (LCS) results are expressed as the percent recovery of analyte. 5. Reporting Quantitation Limits (RQL), typically at or above the Practical Quantitation Limit (PQL) of the analytical method. 6. Method numbers typically denote USEPA procedures. Less than ("<") values reflect nominal quantitation limits adjusted for any required dilutions. 7. Data Qualifiers are J = analyte detected between the RQL and the MDL. B = Analyte detected in associated method blank (s). C=poor CCV recovery. L=poor LCS recovery. S & S1 =MS and/or MSD recovery exceed advisory limits. S2 =Post digestion spike (PDS) recovery exceeds advisory limit. S3 =MS and/or MSD and PDS recoveries exceed advisory limits. P =Precision higher than advisory limit. M =Matrix interference. N=not NELACcertified. N1=subcontract result enquire concerning NELAC certification. Solid sample results for all metals, except Mercury, reported on a dry weight basis (DWB)s. All other results for solid samples reported on an as received basis unless specifically identified as DWB.

Client: Gainco, Inc.
Attn: Paul Supak

Project ID: Falcon Refinery
Sample Name: Tank 20

Report#/Lab ID#: 352492
Sample Matrix: water

REPORT OF ANALYSIS-cont.

QUALITY ASSURANCE DATA ¹

| Parameter | Result | Units | RQL ⁵ | Blank | Date | Method ⁶ | Data Qual. ⁷ | Prec. ² | Recov. ³ | CCV ⁴ | LCS ⁴ |
|--------------------------------------|--------|-------|------------------|-------|----------|---------------------|-------------------------|--------------------|---------------------|------------------|------------------|
| 1,2-Dichlorobenzene | <1 | µg/L | 1 | <1 | 05/13/11 | 8260b & 624 | --- | 2.4 | 111.8 | 110.8 | 112.7 |
| 1,2-Dichloroethane | <1 | µg/L | 1 | <1 | 05/13/11 | 8260b & 624 | --- | 1.5 | 100.6 | 93.6 | 91.9 |
| 1,2-Dichloropropane | <1 | µg/L | 1 | <1 | 05/13/11 | 8260b & 624 | C, | 1.7 | 135.5 | 128.3 | 122.9 |
| 1,3-Dichlorobenzene | <1 | µg/L | 1 | <1 | 05/13/11 | 8260b & 624 | --- | 0.2 | 111.2 | 109.2 | 111.6 |
| 1,4-Dichlorobenzene | <1 | µg/L | 1 | <1 | 05/13/11 | 8260b & 624 | --- | 0.8 | 110.2 | 108.1 | 108.7 |
| 1,4-Dioxane | <20 | µg/L | 20 | <20 | 05/13/11 | 8260b & 624 | --- | 1.9 | 118.4 | 116.9 | 110.4 |
| 2-Butanone (MEK) | 85.4 | µg/L | 20 | <20 | 05/13/11 | 8260b & 624 | --- | 6.9 | 130.2 | 130 | 122.4 |
| 2-Chloroethyl vinyl ether | <1 | µg/L | 1 | <1 | 05/13/11 | 8260b & 624 | --- | 4.5 | 76.6 | 91.1 | 95.4 |
| 2-Hexanone | <2 | µg/L | 2 | <2 | 05/13/11 | 8260b & 624 | --- | 1.7 | 114.7 | 115.7 | 114.6 |
| 4-Methyl-2-pentanone (MIBK) | 6.21 | µg/L | 2 | <2 | 05/13/11 | 8260b & 624 | --- | 0.4 | 125.4 | 123.8 | 118.7 |
| Acetone (2-Propanone) | 624 | µg/L | 10 | <10 | 05/13/11 | 8260b & 624 | --- | 1.9 | 101.8 | 97.1 | 94.9 |
| Acetonitrile | <20 | µg/L | 20 | <20 | 05/13/11 | 8260b & 624 | --- | 0.2 | 126 | 123 | 119.1 |
| Acrolein | <10 | µg/L | 10 | <10 | 05/13/11 | 8260b & 624 | --- | 1 | 88.5 | 112.6 | 100.7 |
| Acrylonitrile | <10 | µg/L | 10 | <10 | 05/13/11 | 8260b & 624 | C, | 0.2 | 141.6 | 140.2 | 131.3 |
| Benzene | 7.78 | µg/L | 1 | <1 | 05/13/11 | 8260b & 624 | --- | 1.8 | 130.2 | 122.1 | 116.9 |
| Bromobenzene | <1 | µg/L | 1 | <1 | 05/13/11 | 8260b & 624 | --- | 0 | 121.1 | 118.6 | 120 |
| Bromodichloromethane | <1 | µg/L | 1 | <1 | 05/13/11 | 8260b & 624 | --- | 2.1 | 108.7 | 101.4 | 101.8 |
| Bromoform (Tribromomethane) | <1 | µg/L | 1 | <1 | 05/13/11 | 8260b & 624 | --- | 4 | 93.6 | 94 | 93.2 |
| Bromomethane (Methyl bromide) | <2 | µg/L | 2 | <2 | 05/13/11 | 8260b & 624 | --- | 10 | 57.9 | 60.8 | 76 |
| Carbon disulfide | <2 | µg/L | 2 | <2 | 05/13/11 | 8260b & 624 | --- | 12.7 | 127.5 | 106 | 112.5 |
| Carbon tetrachloride | <1 | µg/L | 1 | <1 | 05/13/11 | 8260b & 624 | --- | 5.5 | 93.4 | 86.3 | 87.2 |
| Chlorobenzene | <1 | µg/L | 1 | <1 | 05/13/11 | 8260b & 624 | --- | 2.9 | 114.4 | 110.1 | 111 |
| Chloroethane | <2 | µg/L | 2 | <2 | 05/13/11 | 8260b & 624 | --- | 3.8 | 103.9 | 97.9 | 90.4 |
| Chloroform (Trichloromethane) | <1 | µg/L | 1 | <1 | 05/13/11 | 8260b & 624 | --- | 2.5 | 122 | 114.1 | 113.7 |
| Chloromethane (Methyl chloride) | <2 | µg/L | 2 | <2 | 05/13/11 | 8260b & 624 | J, | 5 | 92 | 85.7 | 89.9 |
| cis-1,2-Dichloroethene | <1 | µg/L | 1 | <1 | 05/13/11 | 8260b & 624 | --- | 1.7 | 131.5 | 121.4 | 118.1 |
| cis-1,3-Dichloropropene | <1 | µg/L | 1 | <1 | 05/13/11 | 8260b & 624 | --- | 0.8 | 101.1 | 112.1 | 106.9 |
| Dibromochloromethane | <1 | µg/L | 1 | <1 | 05/13/11 | 8260b & 624 | --- | 0.9 | 106.3 | 104 | 104.3 |
| Dibromomethane (Methylene bromide) | <1 | µg/L | 1 | <1 | 05/13/11 | 8260b & 624 | --- | 2.6 | 113.3 | 106 | 104.7 |
| Dichlorodifluoromethane | <1 | µg/L | 1 | <1 | 05/13/11 | 8260b & 624 | --- | 1.1 | 45.8 | 44 | 44.2 |
| Ethylbenzene | 3.33 | µg/L | 1 | <1 | 05/13/11 | 8260b & 624 | --- | 3.6 | 117.2 | 114.2 | 113.2 |
| Iodomethane (Methyl iodide) | <2 | µg/L | 2 | <2 | 05/13/11 | 8260b & 624 | --- | 1.9 | 57.8 | 60 | 68.9 |
| m,p-Xylenes | 2.41 | µg/L | 2 | <2 | 05/13/11 | 8260b & 624 | --- | 2.7 | 117.1 | 111.6 | 109.6 |
| Methylene chloride (Dichloromethane) | <2 | µg/L | 2 | <2 | 05/13/11 | 8260b & 624 | J, | 16.5 | 131.7 | 112.6 | 112.2 |

Client: Gainco, Inc.
Attn: Paul Supak

Project ID: Falcon Refinery
Sample Name: Tank 20

Report#/Lab ID#: 352492
Sample Matrix: water

REPORT OF ANALYSIS-cont.

QUALITY ASSURANCE DATA ¹

| Parameter | Result | Units | RQL ⁵ | Blank | Date | Method ⁶ | Data Qual. ⁷ | Prec. ² | Recov. ³ | CCV ⁴ | LCS ⁴ |
|---------------------------------------|--------|-------|------------------|-------|----------|---------------------|-------------------------|--------------------|---------------------|------------------|------------------|
| MTBE | 133 | µg/L | 10 | <10 | 05/13/11 | 8260b & 624 | --- | 2.6 | 127 | 119.3 | 111.2 |
| o-Xylene | 1.59 | µg/L | 1 | <1 | 05/13/11 | 8260b & 624 | --- | 1.5 | 118.4 | 111.5 | 113.3 |
| Styrene | 2.29 | µg/L | 1 | <1 | 05/13/11 | 8260b & 624 | --- | 1.2 | 119.4 | 113.9 | 113.3 |
| Tetrachloroethene (Perchloroethylene) | <1 | µg/L | 1 | <1 | 05/13/11 | 8260b & 624 | --- | 4.5 | 104.4 | 100.8 | 100.4 |
| Toluene | 2.79 | µg/L | 1 | <1 | 05/13/11 | 8260b & 624 | --- | 3.5 | 122.3 | 115.7 | 112.7 |
| trans-1,2-Dichloroethene | <1 | µg/L | 1 | <1 | 05/13/11 | 8260b & 624 | --- | 2.4 | 129.2 | 119.7 | 127.3 |
| trans-1,3-Dichloropropene | <1 | µg/L | 1 | <1 | 05/13/11 | 8260b & 624 | --- | 2 | 102.6 | 108.6 | 112.3 |
| Trichloroethene | <1 | µg/L | 1 | <1 | 05/13/11 | 8260b & 624 | --- | 3.3 | 117.7 | 107.9 | 103.5 |
| Trichlorofluoromethane | <1 | µg/L | 1 | <1 | 05/13/11 | 8260b & 624 | --- | 4.5 | 96.1 | 86.5 | 92.6 |
| Vinyl acetate | <2 | µg/L | 2 | <2 | 05/13/11 | 8260b & 624 | --- | 2.2 | 112.2 | 127.4 | 125 |
| Vinyl chloride | <1 | µg/L | 1 | <1 | 05/13/11 | 8260b & 624 | --- | 5.3 | 86.8 | 80.4 | 83.2 |
| Extractable organics-625/8270 | --- | --- | --- | --- | 05/14/11 | 8270c & 625 | --- | --- | --- | --- | --- |
| 1-Methylnaphthalene | <10 | µg/L | 10 | <10 | 05/14/11 | 8270c & 625 | N, | 8 | 28.8 | 88.1 | 45.2 |
| 1,2-Diphenylhydrazine | <10 | µg/L | 10 | <10 | 05/14/11 | 8270c & 625 | --- | 15.7 | 55.8 | 92.5 | 78.6 |
| 1,2,4-Trichlorobenzene | <10 | µg/L | 10 | <10 | 05/14/11 | 8270c & 625 | --- | 3.3 | 21.6 | 90 | 37.6 |
| 2-Chloronaphthalene | <10 | µg/L | 10 | <10 | 05/14/11 | 8270c & 625 | --- | 9 | 29.1 | 95.2 | 46.2 |
| 2-Chlorophenol | <10 | µg/L | 10 | <10 | 05/14/11 | 8270c & 625 | --- | 7.8 | 40.2 | 93.2 | 57.1 |
| 2-Methylnaphthalene | <10 | µg/L | 10 | <10 | 05/14/11 | 8270c & 625 | --- | 9.8 | 27.9 | 90.4 | 43.3 |
| 2-Methylphenol (o-Cresol) | 138 | µg/L | 100 | <100 | 05/14/11 | 8270c & 625 | --- | 22.2 | 46.5 | 93.9 | 63.6 |
| 2-Nitroaniline | <50 | µg/L | 50 | <50 | 05/14/11 | 8270c & 625 | --- | 9.2 | 52.6 | 93.9 | 81.5 |
| 2-Nitrophenol | <10 | µg/L | 10 | <10 | 05/14/11 | 8270c & 625 | --- | 5.9 | 36.6 | 98 | 56.1 |
| 2,4-Dichlorophenol | <10 | µg/L | 10 | <10 | 05/14/11 | 8270c & 625 | --- | 6.5 | 41.4 | 95.9 | 65.9 |
| 2,4-Dimethylphenol | <10 | µg/L | 10 | <10 | 05/14/11 | 8270c & 625 | --- | 3.8 | 48.3 | 86.8 | 78.3 |
| 2,4-Dinitrophenol | <10 | µg/L | 10 | <10 | 05/14/11 | 8270c & 625 | --- | 18.1 | 19.8 | 98.7 | 34.6 |
| 2,4-Dinitrotoluene | <10 | µg/L | 10 | <10 | 05/14/11 | 8270c & 625 | --- | 13.3 | 52.2 | 102.9 | 77.7 |
| 2,4,6-Trichlorophenol | <10 | µg/L | 10 | <10 | 05/14/11 | 8270c & 625 | --- | 10.4 | 39.4 | 96.4 | 56 |
| 2,6-Dichlorophenol | <10 | µg/L | 10 | <10 | 05/14/11 | 8270c & 625 | --- | 3.3 | 36.6 | 92.3 | 52.6 |
| 2,6-Dinitrotoluene | <10 | µg/L | 10 | <10 | 05/14/11 | 8270c & 625 | --- | 16.2 | 55.8 | 97.8 | 82.3 |
| 3-Nitroaniline | <50 | µg/L | 50 | <50 | 05/14/11 | 8270c & 625 | --- | 8.7 | 50.9 | 102 | 82.9 |
| 3,3'-Dichlorobenzidine | <20 | µg/L | 20 | <20 | 05/14/11 | 8270c & 625 | --- | 13.3 | 56.2 | 96.6 | 86.9 |
| 3&4 Methylphenol (m&p-Cresol) | 165 | µg/L | 20 | <20 | 05/14/11 | 8270c & 625 | --- | 15.8 | 40.8 | 96 | 59.6 |
| 4-Bromophenyl phenyl ether | <10 | µg/L | 10 | <10 | 05/14/11 | 8270c & 625 | --- | 19.5 | 52.4 | 95.6 | 70.8 |
| 4-Chloro-3-methylphenol | <20 | µg/L | 20 | <20 | 05/14/11 | 8270c & 625 | --- | 4.1 | 51.6 | 92.8 | 73.9 |

Client: Gainco, Inc.
Attn: Paul Supak

Project ID: Falcon Refinery
Sample Name: Tank 20

Report#/Lab ID#: 352492
Sample Matrix: water

REPORT OF ANALYSIS-cont.
QUALITY ASSURANCE DATA ¹

| Parameter | Result | Units | RQL ⁵ | Blank | Date | Method ⁶ | Data Qual. ⁷ | Prec. ² | Recov. ³ | CCV ⁴ | LCS ⁴ |
|---|--------|-------|------------------|-------|----------|---------------------|-------------------------|--------------------|---------------------|------------------|------------------|
| 4-Chloroaniline (p-Chloroaniline) | <20 | µg/L | 20 | <20 | 05/14/11 | 8270c & 625 | --- | 15.7 | 36.3 | 106.7 | 77.6 |
| 4-Chlorophenyl phenyl ether | <10 | µg/L | 10 | <10 | 05/14/11 | 8270c & 625 | --- | 13.2 | 45.8 | 96.5 | 66.6 |
| 4-Nitroaniline | <20 | µg/L | 20 | <20 | 05/14/11 | 8270c & 625 | --- | 6.7 | 44.8 | 97.5 | 73.3 |
| 4-Nitrophenol | <50 | µg/L | 50 | <50 | 05/14/11 | 8270c & 625 | --- | 7.2 | 12.6 | 83 | 16.2 |
| 4,6-Dinitro-2-methylphenol | <25 | µg/L | 25 | <25 | 05/14/11 | 8270c & 625 | --- | 19.5 | 35.5 | 101.9 | 50.3 |
| 7,12-Dimethylbenz[a]anthracene | <10 | µg/L | 10 | <10 | 05/14/11 | 8270c & 625 | --- | 14.5 | 56.9 | 99.4 | 84.4 |
| Acenaphthene | <10 | µg/L | 10 | <10 | 05/14/11 | 8270c & 625 | --- | 12.8 | 40.6 | 94.3 | 58.1 |
| Acenaphthylene | <10 | µg/L | 10 | <10 | 05/14/11 | 8270c & 625 | --- | 8.4 | 40.9 | 96.1 | 61.8 |
| Aniline | <10 | µg/L | 10 | <10 | 05/14/11 | 8270c & 625 | --- | 15.2 | 47.1 | 92.1 | 59 |
| Anthracene | <10 | µg/L | 10 | <10 | 05/14/11 | 8270c & 625 | --- | 16.1 | 58.8 | 93.4 | 82.7 |
| Benzidine | <40 | µg/L | 40 | <40 | 05/14/11 | 8270c & 625 | --- | 14.9 | 54.8 | 85.5 | 93.8 |
| Benzo[a]anthracene | <10 | µg/L | 10 | <10 | 05/14/11 | 8270c & 625 | J, | 11 | 56.4 | 93.7 | 84.6 |
| Benzo[a]pyrene | <10 | µg/L | 10 | <10 | 05/14/11 | 8270c & 625 | J, | 10.9 | 57.9 | 97 | 86 |
| Benzo[b]fluoranthene | <10 | µg/L | 10 | <10 | 05/14/11 | 8270c & 625 | --- | 13.2 | 55 | 97.6 | 81.8 |
| Benzo[g,h,i]perylene | <10 | µg/L | 10 | <10 | 05/14/11 | 8270c & 625 | --- | 12.2 | 64.6 | 90.3 | 101.1 |
| Benzo[j,k]fluoranthene | <10 | µg/L | 10 | <10 | 05/14/11 | 8270c & 625 | J, | 13.9 | 56.2 | 94.7 | 83.4 |
| Benzoic acid | <40 | µg/L | 40 | <40 | 05/14/11 | 8270c & 625 | --- | 26.7 | 1.9 | 88.3 | 4 |
| Benzyl alcohol | <20 | µg/L | 20 | <20 | 05/14/11 | 8270c & 625 | --- | 19.9 | 38 | 96.5 | 71.7 |
| bis(2-Chloroethoxy)methane | <10 | µg/L | 10 | <10 | 05/14/11 | 8270c & 625 | --- | 5 | 49.2 | 90.9 | 80.7 |
| bis(2-Chloroethyl)ether | <10 | µg/L | 10 | <10 | 05/14/11 | 8270c & 625 | --- | 6.5 | 44.7 | 93.2 | 66.3 |
| bis(2-chloroisopropyl)ether | <10 | µg/L | 10 | <10 | 05/14/11 | 8270c & 625 | --- | 9.8 | 40.2 | 90.9 | 61.4 |
| bis(2-Ethylhexyl)phthalate | <10 | µg/L | 10 | <10 | 05/14/11 | 8270c & 625 | --- | 14.1 | 53.2 | 98.3 | 83.1 |
| Butyl benzyl phthalate | <10 | µg/L | 10 | <10 | 05/14/11 | 8270c & 625 | --- | 6.9 | 50.9 | 95.3 | 83.8 |
| Chrysene | 17.2 | µg/L | 10 | <10 | 05/14/11 | 8270c & 625 | --- | 9.5 | 55.6 | 91.8 | 86.1 |
| Di-n-butyl phthalate (Dibutylphthalate) | <10 | µg/L | 10 | <10 | 05/14/11 | 8270c & 625 | --- | 15.5 | 54.3 | 91 | 80.4 |
| Di-n-octylphthalate (Dioctylphthalate) | <10 | µg/L | 10 | <10 | 05/14/11 | 8270c & 625 | --- | 11.9 | 52.7 | 101.8 | 82.1 |
| Dibenz[a,h]acridine | <10 | µg/L | 10 | <10 | 05/14/11 | 8270c & 625 | --- | 16.6 | 59.1 | 93.3 | 89.1 |
| Dibenz[a,h]anthracene | <10 | µg/L | 10 | <10 | 05/14/11 | 8270c & 625 | --- | 13.3 | 64.2 | 90.5 | 95.7 |
| Dibenzofuran | <10 | µg/L | 10 | <10 | 05/14/11 | 8270c & 625 | --- | 13 | 44.2 | 96 | 65.2 |
| Diethylphthalate | <10 | µg/L | 10 | <10 | 05/14/11 | 8270c & 625 | --- | 15 | 53.6 | 96.9 | 75.5 |
| Dimethylphthalate | <10 | µg/L | 10 | <10 | 05/14/11 | 8270c & 625 | --- | 9.2 | 53 | 94.9 | 77.6 |
| Fluoranthene | <10 | µg/L | 10 | <10 | 05/14/11 | 8270c & 625 | J, | 11.8 | 57.5 | 95.1 | 84.3 |
| Fluorene | <10 | µg/L | 10 | <10 | 05/14/11 | 8270c & 625 | --- | 9.2 | 49.9 | 97.1 | 72.3 |
| Hexachlorobenzene | <10 | µg/L | 10 | <10 | 05/14/11 | 8270c & 625 | --- | 12.4 | 55.9 | 93.6 | 75.6 |

Client: Gainco, Inc.
Attn: Paul Supak

Project ID: Falcon Refinery
Sample Name: Tank 20

Report#/Lab ID#: 352492
Sample Matrix: water

REPORT OF ANALYSIS-cont.
QUALITY ASSURANCE DATA ¹

| Parameter | Result | Units | RQL ⁵ | Blank | Date | Method ⁶ | Data Qual. ⁷ | Prec. ² | Recov. ³ | CCV ⁴ | LCS ⁴ |
|-----------------------------------|--------|-------|------------------|-------|----------|---------------------|-------------------------|--------------------|---------------------|------------------|------------------|
| Hexachlorobutadiene | <10 | µg/L | 10 | <10 | 05/14/11 | 8270c & 625 | --- | 1.9 | 19.5 | 93 | 31.8 |
| Hexachlorocyclopentadiene (HCCPD) | <10 | µg/L | 10 | <10 | 05/14/11 | 8270c & 625 | S1,M, | 0 | 0 | 102.8 | 14.8 |
| Hexachloroethane | <10 | µg/L | 10 | <10 | 05/14/11 | 8270c & 625 | --- | 2.9 | 19.7 | 94.8 | 31.4 |
| Indene | <10 | µg/L | 10 | <10 | 05/14/11 | 8270c & 625 | N, | 11.5 | 34.2 | 93.3 | 50.2 |
| Indeno[1,2,3-cd]pyrene | <10 | µg/L | 10 | <10 | 05/14/11 | 8270c & 625 | --- | 16.1 | 61 | 92.2 | 89.4 |
| Isophorone | <10 | µg/L | 10 | <10 | 05/14/11 | 8270c & 625 | --- | 6.7 | 57.6 | 90.4 | 95.3 |
| Methylchrysene | <10 | µg/L | 10 | <10 | 05/14/11 | 8270c & 625 | J,N, | 10.2 | 46.9 | 91.9 | 69.2 |
| N-Nitrosodi-n-propylamine | <10 | µg/L | 10 | <10 | 05/14/11 | 8270c & 625 | --- | 4.4 | 43.1 | 88.7 | 71.9 |
| N-Nitrosodimethylamine | <10 | µg/L | 10 | <10 | 05/14/11 | 8270c & 625 | --- | 5.9 | 31 | 94.3 | 41.5 |
| N-Nitrosodiphenylamine | <10 | µg/L | 10 | <10 | 05/14/11 | 8270c & 625 | --- | 14.4 | 59.5 | 91.7 | 85.4 |
| Naphthalene | <10 | µg/L | 10 | <10 | 05/14/11 | 8270c & 625 | J, | 5.5 | 31.7 | 90.7 | 49.8 |
| Nitrobenzene | <10 | µg/L | 10 | <10 | 05/14/11 | 8270c & 625 | --- | 1.6 | 44.1 | 91.2 | 72.3 |
| Pentachlorophenol | <10 | µg/L | 10 | <10 | 05/14/11 | 8270c & 625 | --- | 9.5 | 33.2 | 101.3 | 54.8 |
| Phenanthrene | 14.7 | µg/L | 10 | <10 | 05/14/11 | 8270c & 625 | --- | 12.2 | 59.8 | 92.6 | 86 |
| Phenol | 509 | µg/L | 100 | <100 | 05/14/11 | 8270c & 625 | --- | 15.4 | 17.9 | 93.1 | 25 |
| Pyrene | 18.7 | µg/L | 10 | <10 | 05/14/11 | 8270c & 625 | --- | 11.1 | 58.7 | 95.5 | 85.8 |
| Pyridine | <10 | µg/L | 10 | <10 | 05/14/11 | 8270c & 625 | --- | 19.5 | 26.1 | 94.6 | 34.1 |
| Quinoline | <10 | µg/L | 10 | <10 | 05/14/11 | 8270c & 625 | --- | 9.8 | 51.2 | 91.8 | 79.4 |

Client: Gainco, Inc.
Attn: Paul Supak

Project ID: Falcon Refinery
Sample Name: Tank 20

Report#/Lab ID#: 352492
Sample Matrix: water

REPORT OF SURROGATE RECOVERY

| Surrogate Compound | Method | Recovery | Recovery Limits | Date Analyzed | Data Qualifiers |
|-----------------------|-------------|--------------|-----------------|---------------|-----------------|
| 1,2-Dichloroethane-d4 | 8260b & 624 | 85.4 | 70-125 | 05/13/11 | --- |
| 1,2-Dichloroethane-d4 | 8260b & 624 | 81.5 | 70-125 | 05/13/11 | --- |
| 4-Bromofluorobenzene | 8260b & 624 | 100.6 | 80-115 | 05/13/11 | --- |
| 4-Bromofluorobenzene | 8260b & 624 | 94.8 | 80-115 | 05/13/11 | --- |
| Toluene-d8 | 8260b & 624 | 103.9 | 78-115 | 05/13/11 | --- |
| Toluene-d8 | 8260b & 624 | 105.3 | 78-115 | 05/13/11 | --- |
| 2-Fluorobiphenyl | 8270c & 625 | 40 | 10-110 | 05/14/11 | --- |
| 2-Fluorobiphenyl | 8270c & 625 | none/diluted | diluted @ 10X | 05/14/11 | D |
| 2-Fluorophenol | 8270c & 625 | 28.2 | 10-110 | 05/14/11 | --- |
| 2-Fluorophenol | 8270c & 625 | none/diluted | diluted @ 10X | 05/14/11 | D |
| 2,4,6-Tribromophenol | 8270c & 625 | 51.2 | 10-120 | 05/14/11 | --- |
| 2,4,6-Tribromophenol | 8270c & 625 | none/diluted | diluted @ 10X | 05/14/11 | D |
| Nitrobenzene-d5 | 8270c & 625 | 45.6 | 10-110 | 05/14/11 | --- |
| Nitrobenzene-d5 | 8270c & 625 | none/diluted | diluted @ 10X | 05/14/11 | D |
| Phenol-d6 | 8270c & 625 | 23 | 10-110 | 05/14/11 | --- |
| Phenol-d6 | 8270c & 625 | none/diluted | diluted @ 10X | 05/14/11 | D |
| Terphenyl-d14 | 8270c & 625 | 27.3 | 10-115 | 05/14/11 | --- |
| Terphenyl-d14 | 8270c & 625 | none/diluted | diluted @ 10X | 05/14/11 | D |

Data Qualifiers: D= Surrogates diluted and X= Surrogates outside advisory recovery limits.

Exceptions Report (FINAL SECTION / END-OF-REPORT):**Report #/Lab ID#:** 352492 **Matrix:** water**Client:** Gainco, Inc.**Attn:** Paul Supak**Project ID:** Falcon Refinery**Sample Name:** Tank 20

Unless otherwise identified by data qualifier "N" or by an exception report, all reported results represent parameters and tests for which AnalySys maintains NELAC certification; or results provided by a subcontractor with NELAC certification for the test results provided.

**Sample Temperature/Condition:** ≤6°C

The typical sample temperature criteria (except for metals by ICP, GFAA and AA and a very few other tests) is ≤ 6°C. Possible exceptions include samples submitted to laboratory within such a short time after sampling that cooling measures used in the field and during transport had insufficient time to achieve desired temperatures in the samples (see sample collection and sample receipt times) and samples where the temperature could not be measured due to sample submission in a manner precluding temperature measurement without impacting sample integrity (ex. in a bottle with no cooler).

Standard sample acceptability conditions met? : YES

Sample received in appropriate container(s), at appropriate temperature and pH.

J flag Discussion:

A J-flag data qualifier indicates that the raw calculated analyte concentration in the sample (uncorrected for background levels/blanks and other potential sources of sampling and analytical contamination), though less than the Reported Quantitation Limit (RQL) is greater than the Detection Limit. Because the reported result is below the quantitation limit for this project/sample (or test procedure), GC/MS organics results may or MAY NOT have been verified as to the presence and relative ratio of target ions (eg. the material causing the J flag "hit" in such situations may be nothing more than background ion-fragment noise.)

Comments pertaining to Data Qualifiers and QC data (where applicable):

| Parameter | Qualif. | Comments |
|--------------------------------------|---------|---|
| Chromium/ICPMS | J | See J-flag discussion above. |
| Selenium/ICPMS | J | See J-flag discussion above. |
| 1,1,1-Trichloroethane | L | Lab control sample (LCS or spiked blank). LCS recov-high (high bias). Sample result < MDL. No impact. |
| 1,1,1-Trichloroethane | S | Spike (MS,MSD,PDS) recovery issue. MS, MSD & PDS recovery outside acceptance range. LCS fails or not available. Probable sample impact. |
| 1,1,2,2-Tetrachloroethane | C | Cont. Calib. Verification (CCV). CCV recov-high (high bias). Sample result < MDL. No impact. |
| 1,2-Dichloropropane | C | Cont. Calib. Verification (CCV). CCV recov-high (high bias). Sample result < MDL. No impact. |
| Acrylonitrile | C | Cont. Calib. Verification (CCV). CCV recov-high (high bias). Sample result < MDL. No impact. |
| Chloromethane (Methyl chloride) | J | See J-flag discussion above. |
| Methylene chloride (Dichloromethane) | J | See J-flag discussion above. |
| 1-Methylnaphthalene | N | NELAC accreditation for this analyte not available from TCEQ. 30 TAC§25.6(4) applies. |
| Benzo[a]anthracene | J | See J-flag discussion above. |
| Benzo[a]pyrene | J | See J-flag discussion above. |
| Benzo[j,k]fluoranthene | J | See J-flag discussion above. |
| Fluoranthene | J | See J-flag discussion above. |
| Hexachlorocyclopentadiene (HCCPD) | S1 | Spike (MS,MSD) recovery issue. MS & MSD recovery outside acceptance range. LCS recovery OK. Probable matrix interference. |
| Indene | N | NELAC accreditation for this analyte not available from TCEQ. 30 TAC§25.6(4) applies. |
| Methylchrysene | J | See J-flag discussion above. |
| Methylchrysene | N | NELAC accreditation for this analyte not available from TCEQ. 30 TAC§25.6(4) applies. |
| Naphthalene | J | See J-flag discussion above. |

Exceptions Report (FINAL SECTION / END-OF-REPORT):

Report #/Lab ID#: 352492 Matrix: water

Client: Gainco, Inc.

Attn: Paul Supak

Project ID: Falcon Refinery

Sample Name: Tank 20

Unless otherwise identified by data qualifier "N" or by an exception report, all reported results represent parameters and tests for which AnalySys maintains NELAC certification; or results provided by a subcontractor with NELAC certification for the test results provided.



| | | |
|----------------------|---|---|
| 2-Fluorobiphenyl | D | Surrogate recoveries not accurately quantifiable. |
| 2-Fluorophenol | D | Surrogate recoveries not accurately quantifiable. |
| 2,4,6-Tribromophenol | D | Surrogate recoveries not accurately quantifiable. |
| Nitrobenzene-d5 | D | Surrogate recoveries not accurately quantifiable. |
| Phenol-d6 | D | Surrogate recoveries not accurately quantifiable. |
| Terphenyl-d14 | D | Surrogate recoveries not accurately quantifiable. |

Client: Gainco, Inc.
Attn: Paul Supak
Address:

Phone: **FAX:**

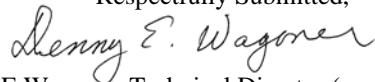
Report#/Lab ID#: 352493 **Report Date:** 05/16/11
Project ID: Falcon Refinery
Sample Name: Tank 2
Sample Matrix: water
Date Received: 05/10/2011 **Time:** 16:11
Date Sampled: 05/10/2011 **Time:** 11:15

REPORT OF ANALYSIS

QUALITY ASSURANCE DATA ¹

| Parameter | Result | Units | RQL ⁵ | Blank | Date | Method ⁶ | Data Qual. ⁷ | Prec. ² | Recov. ³ | CCV ⁴ | LCS ⁴ |
|--|---------|-------|------------------|---------|----------|---------------------|-------------------------|--------------------|---------------------|------------------|------------------|
| A/BN extraction-625/8270 | --- | --- | --- | --- | 05/13/11 | 3510 | --- | --- | --- | --- | --- |
| Metals Dig.-Hg | --- | --- | --- | --- | 05/11/11 | 7470 & SM3112B | --- | --- | --- | --- | --- |
| Metals Dig.-Total | --- | --- | --- | --- | 05/11/11 | 200.2 & 3005A | --- | --- | --- | --- | --- |
| Arsenic/ICPMS | 0.00338 | mg/L | 0.002 | <0.002 | 05/13/11 | 6020A & 200.8 | --- | 0 | 99.5 | 98.3 | 99.6 |
| Barium/ICPMS | 0.474 | mg/L | 0.002 | <0.002 | 05/13/11 | 6020A & 200.8 | --- | 1 | 104.9 | 106.6 | 100 |
| Cadmium/ICPMS | <0.001 | mg/L | 0.001 | <0.001 | 05/13/11 | 6020A & 200.8 | --- | 0.8 | 93.9 | 100.5 | 92.7 |
| Chromium/ICPMS | 0.00491 | mg/L | 0.002 | <0.002 | 05/13/11 | 6020A & 200.8 | --- | 1.5 | 99.5 | 98.5 | 100 |
| Lead/ICPMS | 0.0252 | mg/L | 0.002 | <0.002 | 05/13/11 | 6020A & 200.8 | --- | 0.1 | 87 | 95.6 | 92.2 |
| Mercury/CVAA | <0.0002 | mg/L | 0.0002 | <0.0002 | 05/12/11 | 7470&SM3112B | --- | 3.08 | 90 | 107 | 107 |
| Selenium/ICPMS | <0.002 | mg/L | 0.002 | <0.002 | 05/13/11 | 6020A & 200.8 | --- | 1.7 | 106.2 | 104 | 106.9 |
| Silver/ICPMS | <0.001 | mg/L | 0.001 | <0.001 | 05/13/11 | 6020A & 200.8 | --- | 1 | 92.7 | 97.6 | 92.2 |
| Volatile organics-8260 | --- | --- | --- | --- | 05/12/11 | 8260b(5030/5035) | --- | --- | --- | --- | --- |
| 1,1-Dichloroethane | <10 | µg/L | 10 | <10 | 05/12/11 | 8260b & 624 | --- | 5.9 | 136.5 | 119.2 | 121.3 |
| 1,1-Dichloroethene | <10 | µg/L | 10 | <10 | 05/12/11 | 8260b & 624 | --- | 7.3 | 107.9 | 94.5 | 100.3 |
| 1,1,1-Trichloroethane | <10 | µg/L | 10 | <10 | 05/12/11 | 8260b & 624 | C,L,S, | 8.2 | 143.9 | 127.9 | 125.8 |
| 1,1,1,2-Tetrachloroethane | <10 | µg/L | 10 | <10 | 05/12/11 | 8260b & 624 | --- | 1 | 107.3 | 101.8 | 102.7 |
| 1,1,2-Trichloroethane | <10 | µg/L | 10 | <10 | 05/12/11 | 8260b & 624 | --- | 2.2 | 124.7 | 116.1 | 114.5 |
| 1,1,2,2-Tetrachloroethane | <10 | µg/L | 10 | <10 | 05/12/11 | 8260b & 624 | --- | 1.6 | 129.1 | 128.8 | 122.8 |
| 1,2-Dibromo-3-chloropropane | <20 | µg/L | 20 | <20 | 05/12/11 | 8260b & 624 | --- | 2.6 | 112.3 | 104.9 | 106.1 |
| 1,2-Dibromoethane (Ethylene dibromide) | <10 | µg/L | 10 | <10 | 05/12/11 | 8260b & 624 | --- | 0.5 | 118.1 | 112 | 112.5 |
| 1,2-Dichlorobenzene | <10 | µg/L | 10 | <10 | 05/12/11 | 8260b & 624 | --- | 1.2 | 111.4 | 110.8 | 110 |

This analytical report is respectfully submitted by AnalySys, Inc. The enclosed results reflect only the sample identified above. The results have been carefully reviewed and to the best of my knowledge, unless otherwise indicated, meet NELAC requirements as described by AnalySys, Inc.'s Quality Assurance/Quality Control Program. © Copyright 2003, AnalySys, Inc., Austin, TX. All rights reserved. No part of this publication may be reproduced or transmitted in any form or by any means without the express written consent of AnalySys, Inc.

Respectfully Submitted,

D.E. Wagoner, Technical Director (or designee)

1. Quality assurance data for the sample batch which included this sample. 2. Precision (PREC) is the absolute value of the relative percent difference between duplicate results. 3. Recovery (Recov.) is the percent of analyte recovered from a spiked sample. 4. Calibration Verification (CCV) and Laboratory Control Sample (LCS) results are expressed as the percent recovery of analyte. 5. Reporting Quantitation Limits (RQL), typically at or above the Practical Quantitation Limit (PQL) of the analytical method. 6. Method numbers typically denote USEPA procedures. Less than ("<") values reflect nominal quantitation limits adjusted for any required dilutions. 7. Data Qualifiers are J = analyte detected between the RQL and the MDL. B = Analyte detected in associated method blank (s). C=poor CCV recovery. L=poor LCS recovery. S & S1 =MS and/or MSD recovery exceed advisory limits. S2 =Post digestion spike (PDS) recovery exceeds advisory limit. S3 =MS and/or MSD and PDS recoveries exceed advisory limits. P =Precision higher than advisory limit. M =Matrix interference. N=not NELACcertified. N1=subcontract result enquire concerning NELAC certification. Solid sample results for all metals, except Mercury, reported on a dry weight basis (DWB)s. All other results for solid samples reported on an as received basis unless specifically identified as DWB.

Client: Gainco, Inc.
Attn: Paul Supak

Project ID: Falcon Refinery
Sample Name: Tank 2

Report#/Lab ID#: 352493
Sample Matrix: water

REPORT OF ANALYSIS-cont.
QUALITY ASSURANCE DATA ¹

| Parameter | Result | Units | RQL ⁵ | Blank | Date | Method ⁶ | Data Qual. ⁷ | Prec. ² | Recov. ³ | CCV ⁴ | LCS ⁴ |
|--------------------------------------|--------|-------|------------------|-------|----------|---------------------|-------------------------|--------------------|---------------------|------------------|------------------|
| 1,2-Dichloroethane | <10 | µg/L | 10 | <10 | 05/12/11 | 8260b & 624 | --- | 0.8 | 103.3 | 93.3 | 90.9 |
| 1,2-Dichloropropane | <10 | µg/L | 10 | <10 | 05/12/11 | 8260b & 624 | C, | 2 | 132.1 | 120.6 | 117.1 |
| 1,3-Dichlorobenzene | <10 | µg/L | 10 | <10 | 05/12/11 | 8260b & 624 | --- | 0.1 | 109.7 | 108.2 | 107.9 |
| 1,4-Dichlorobenzene | <10 | µg/L | 10 | <10 | 05/12/11 | 8260b & 624 | --- | 1 | 108.5 | 106.9 | 105.3 |
| 1,4-Dioxane | <200 | µg/L | 200 | <200 | 05/12/11 | 8260b & 624 | --- | 2.5 | 119.4 | 105 | 118.4 |
| 2-Butanone (MEK) | <20 | µg/L | 20 | <20 | 05/12/11 | 8260b & 624 | --- | 2.6 | 137.8 | 115.3 | 128.2 |
| 2-Chloroethyl vinyl ether | <10 | µg/L | 10 | <10 | 05/12/11 | 8260b & 624 | --- | 16.2 | 108.9 | 93.4 | 100.4 |
| 2-Hexanone | <20 | µg/L | 20 | <20 | 05/12/11 | 8260b & 624 | --- | 0.6 | 115.7 | 111.4 | 111.7 |
| 4-Methyl-2-pentanone (MIBK) | <20 | µg/L | 20 | <20 | 05/12/11 | 8260b & 624 | --- | 0.8 | 129.4 | 114.8 | 119.6 |
| Acetone (2-Propanone) | <100 | µg/L | 100 | <100 | 05/12/11 | 8260b & 624 | --- | 5.5 | 105.3 | 92.3 | 96 |
| Acetonitrile | <200 | µg/L | 200 | <200 | 05/12/11 | 8260b & 624 | --- | 1.2 | 133.3 | 112.4 | 118.1 |
| Acrolein | <100 | µg/L | 100 | <100 | 05/12/11 | 8260b & 624 | --- | 3.9 | 123 | 106.5 | 107.7 |
| Acrylonitrile | <100 | µg/L | 100 | <100 | 05/12/11 | 8260b & 624 | --- | 0.4 | 139.8 | 122.9 | 123.3 |
| Benzene | 199 | µg/L | 10 | <10 | 05/12/11 | 8260b & 624 | --- | 4.5 | 126.1 | 114.9 | 111.1 |
| Bromobenzene | <10 | µg/L | 10 | <10 | 05/12/11 | 8260b & 624 | --- | 1 | 117.2 | 116 | 111.7 |
| Bromodichloromethane | <10 | µg/L | 10 | <10 | 05/12/11 | 8260b & 624 | --- | 2 | 110.1 | 99.8 | 98.2 |
| Bromoform (Tribromomethane) | <10 | µg/L | 10 | <10 | 05/12/11 | 8260b & 624 | --- | 0.9 | 100.9 | 95.4 | 95.4 |
| Bromomethane (Methyl bromide) | <20 | µg/L | 20 | <20 | 05/12/11 | 8260b & 624 | --- | 4.7 | 91.6 | 80.7 | 86.8 |
| Carbon disulfide | <20 | µg/L | 20 | <20 | 05/12/11 | 8260b & 624 | --- | 7.9 | 127.6 | 113.6 | 128.7 |
| Carbon tetrachloride | <10 | µg/L | 10 | <10 | 05/12/11 | 8260b & 624 | --- | 5 | 97.6 | 89.7 | 89.7 |
| Chlorobenzene | <10 | µg/L | 10 | <10 | 05/12/11 | 8260b & 624 | --- | 1 | 113.7 | 107.7 | 108.4 |
| Chloroethane | <20 | µg/L | 20 | <20 | 05/12/11 | 8260b & 624 | --- | 8.1 | 106.7 | 96.2 | 92.8 |
| Chloroform (Trichloromethane) | <10 | µg/L | 10 | <10 | 05/12/11 | 8260b & 624 | --- | 2.2 | 120.9 | 107.8 | 110 |
| Chloromethane (Methyl chloride) | <20 | µg/L | 20 | <20 | 05/12/11 | 8260b & 624 | --- | 7.3 | 102.7 | 92 | 94.3 |
| cis-1,2-Dichloroethene | <10 | µg/L | 10 | <10 | 05/12/11 | 8260b & 624 | --- | 2.6 | 130.4 | 116.9 | 113.9 |
| cis-1,3-Dichloropropene | <10 | µg/L | 10 | <10 | 05/12/11 | 8260b & 624 | --- | 0.8 | 117.6 | 108 | 104.8 |
| Dibromochloromethane | <10 | µg/L | 10 | <10 | 05/12/11 | 8260b & 624 | --- | 2.7 | 107.5 | 104.3 | 103.9 |
| Dibromomethane (Methylene bromide) | <10 | µg/L | 10 | <10 | 05/12/11 | 8260b & 624 | --- | 2.3 | 116.5 | 103.4 | 101.6 |
| Dichlorodifluoromethane | <10 | µg/L | 10 | <10 | 05/12/11 | 8260b & 624 | --- | 11.8 | 52.2 | 48.2 | 45.7 |
| Ethylbenzene | 166 | µg/L | 10 | <10 | 05/12/11 | 8260b & 624 | --- | 1 | 114.6 | 109.3 | 109.2 |
| Iodomethane (Methyl iodide) | <20 | µg/L | 20 | <20 | 05/12/11 | 8260b & 624 | --- | 4.8 | 86.2 | 76 | 84.2 |
| m,p-Xylenes | 432 | µg/L | 20 | <20 | 05/12/11 | 8260b & 624 | --- | 0.9 | 113 | 107.6 | 107.6 |
| Methylene chloride (Dichloromethane) | <20 | µg/L | 20 | <20 | 05/12/11 | 8260b & 624 | J, | 2.5 | 137.5 | 113.8 | 116.4 |
| MTBE | 290 | µg/L | 10 | <10 | 05/12/11 | 8260b & 624 | --- | 4.2 | 135.2 | 118.3 | 113.3 |

Client: Gainco, Inc.
Attn: Paul Supak

Project ID: Falcon Refinery
Sample Name: Tank 2

Report#/Lab ID#: 352493
Sample Matrix: water

REPORT OF ANALYSIS-cont.

QUALITY ASSURANCE DATA ¹

| Parameter | Result | Units | RQL ⁵ | Blank | Date | Method ⁶ | Data Qual. ⁷ | Prec. ² | Recov. ³ | CCV ⁴ | LCS ⁴ |
|---------------------------------------|--------|-------|------------------|-------|----------|---------------------|-------------------------|--------------------|---------------------|------------------|------------------|
| o-Xylene | 30.5 | µg/L | 10 | <10 | 05/12/11 | 8260b & 624 | --- | 0.7 | 114.6 | 107.5 | 109.5 |
| Styrene | <10 | µg/L | 10 | <10 | 05/12/11 | 8260b & 624 | --- | 1.2 | 116.1 | 111 | 110.4 |
| Tetrachloroethene (Perchloroethylene) | <10 | µg/L | 10 | <10 | 05/12/11 | 8260b & 624 | --- | 4.1 | 105.6 | 99.4 | 100.1 |
| Toluene | 26.3 | µg/L | 10 | <10 | 05/12/11 | 8260b & 624 | --- | 3.7 | 120.5 | 108.7 | 107.4 |
| trans-1,2-Dichloroethene | <10 | µg/L | 10 | <10 | 05/12/11 | 8260b & 624 | --- | 5.3 | 127 | 113.8 | 123.9 |
| trans-1,3-Dichloropropene | <10 | µg/L | 10 | <10 | 05/12/11 | 8260b & 624 | --- | 1.7 | 112.2 | 108.7 | 112.2 |
| Trichloroethene | <10 | µg/L | 10 | <10 | 05/12/11 | 8260b & 624 | --- | 3.1 | 111.7 | 104 | 100.4 |
| Trichlorofluoromethane | <10 | µg/L | 10 | <10 | 05/12/11 | 8260b & 624 | --- | 7.1 | 99.7 | 87.2 | 93.3 |
| Vinyl acetate | <20 | µg/L | 20 | <20 | 05/12/11 | 8260b & 624 | --- | 4.3 | 140.5 | 119.7 | 120.1 |
| Vinyl chloride | <10 | µg/L | 10 | <10 | 05/12/11 | 8260b & 624 | --- | 11 | 95.8 | 85.9 | 84.6 |
| Extractable organics-625/8270 | --- | --- | --- | --- | 05/13/11 | 8270c & 625 | --- | --- | --- | --- | --- |
| 1-Methylnaphthalene | 10.9 | µg/L | 10 | <10 | 05/13/11 | 8270c & 625 | N, | 8 | 28.8 | 88.1 | 45.2 |
| 1,2-Diphenylhydrazine | <10 | µg/L | 10 | <10 | 05/13/11 | 8270c & 625 | --- | 15.7 | 55.8 | 92.5 | 78.6 |
| 1,2,4-Trichlorobenzene | <10 | µg/L | 10 | <10 | 05/13/11 | 8270c & 625 | --- | 3.3 | 21.6 | 90 | 37.6 |
| 2-Chloronaphthalene | <10 | µg/L | 10 | <10 | 05/13/11 | 8270c & 625 | --- | 9 | 29.1 | 95.2 | 46.2 |
| 2-Chlorophenol | <10 | µg/L | 10 | <10 | 05/13/11 | 8270c & 625 | --- | 7.8 | 40.2 | 93.2 | 57.1 |
| 2-Methylnaphthalene | <10 | µg/L | 10 | <10 | 05/13/11 | 8270c & 625 | J, | 9.8 | 27.9 | 90.4 | 43.3 |
| 2-Methylphenol (o-Cresol) | <10 | µg/L | 10 | <10 | 05/13/11 | 8270c & 625 | J, | 22.2 | 46.5 | 93.9 | 63.6 |
| 2-Nitroaniline | <50 | µg/L | 50 | <50 | 05/13/11 | 8270c & 625 | --- | 9.2 | 52.6 | 93.9 | 81.5 |
| 2-Nitrophenol | <10 | µg/L | 10 | <10 | 05/13/11 | 8270c & 625 | --- | 5.9 | 36.6 | 98 | 56.1 |
| 2,4-Dichlorophenol | <10 | µg/L | 10 | <10 | 05/13/11 | 8270c & 625 | --- | 6.5 | 41.4 | 95.9 | 65.9 |
| 2,4-Dimethylphenol | 20.6 | µg/L | 10 | <10 | 05/13/11 | 8270c & 625 | --- | 3.8 | 48.3 | 86.8 | 78.3 |
| 2,4-Dinitrophenol | <10 | µg/L | 10 | <10 | 05/13/11 | 8270c & 625 | --- | 18.1 | 19.8 | 98.7 | 34.6 |
| 2,4-Dinitrotoluene | <10 | µg/L | 10 | <10 | 05/13/11 | 8270c & 625 | --- | 13.3 | 52.2 | 102.9 | 77.7 |
| 2,4,6-Trichlorophenol | <10 | µg/L | 10 | <10 | 05/13/11 | 8270c & 625 | --- | 10.4 | 39.4 | 96.4 | 56 |
| 2,6-Dichlorophenol | <10 | µg/L | 10 | <10 | 05/13/11 | 8270c & 625 | --- | 3.3 | 36.6 | 92.3 | 52.6 |
| 2,6-Dinitrotoluene | <10 | µg/L | 10 | <10 | 05/13/11 | 8270c & 625 | --- | 16.2 | 55.8 | 97.8 | 82.3 |
| 3-Nitroaniline | <50 | µg/L | 50 | <50 | 05/13/11 | 8270c & 625 | --- | 8.7 | 50.9 | 102 | 82.9 |
| 3,3'-Dichlorobenzidine | <20 | µg/L | 20 | <20 | 05/13/11 | 8270c & 625 | --- | 13.3 | 56.2 | 96.6 | 86.9 |
| 3&4 Methylphenol (m&p-Cresol) | <20 | µg/L | 20 | <20 | 05/13/11 | 8270c & 625 | --- | 15.8 | 40.8 | 96 | 59.6 |
| 4-Bromophenyl phenyl ether | <10 | µg/L | 10 | <10 | 05/13/11 | 8270c & 625 | --- | 19.5 | 52.4 | 95.6 | 70.8 |
| 4-Chloro-3-methylphenol | <20 | µg/L | 20 | <20 | 05/13/11 | 8270c & 625 | --- | 4.1 | 51.6 | 92.8 | 73.9 |
| 4-Chloroaniline (p-Chloroaniline) | <20 | µg/L | 20 | <20 | 05/13/11 | 8270c & 625 | --- | 15.7 | 36.3 | 106.7 | 77.6 |

Client: Gainco, Inc.
Attn: Paul Supak

Project ID: Falcon Refinery
Sample Name: Tank 2

Report#/Lab ID#: 352493
Sample Matrix: water

REPORT OF ANALYSIS-cont.
QUALITY ASSURANCE DATA ¹

| Parameter | Result | Units | RQL ⁵ | Blank | Date | Method ⁶ | Data Qual. ⁷ | Prec. ² | Recov. ³ | CCV ⁴ | LCS ⁴ |
|---|--------|-------|------------------|-------|----------|---------------------|-------------------------|--------------------|---------------------|------------------|------------------|
| 4-Chlorophenyl phenyl ether | <10 | µg/L | 10 | <10 | 05/13/11 | 8270c & 625 | --- | 13.2 | 45.8 | 96.5 | 66.6 |
| 4-Nitroaniline | <20 | µg/L | 20 | <20 | 05/13/11 | 8270c & 625 | --- | 6.7 | 44.8 | 97.5 | 73.3 |
| 4-Nitrophenol | <50 | µg/L | 50 | <50 | 05/13/11 | 8270c & 625 | --- | 7.2 | 12.6 | 83 | 16.2 |
| 4,6-Dinitro-2-methylphenol | <25 | µg/L | 25 | <25 | 05/13/11 | 8270c & 625 | --- | 19.5 | 35.5 | 101.9 | 50.3 |
| 7,12-Dimethylbenz[a]anthracene | <10 | µg/L | 10 | <10 | 05/13/11 | 8270c & 625 | --- | 14.5 | 56.9 | 99.4 | 84.4 |
| Acenaphthene | <10 | µg/L | 10 | <10 | 05/13/11 | 8270c & 625 | --- | 12.8 | 40.6 | 94.3 | 58.1 |
| Acenaphthylene | <10 | µg/L | 10 | <10 | 05/13/11 | 8270c & 625 | --- | 8.4 | 40.9 | 96.1 | 61.8 |
| Aniline | <10 | µg/L | 10 | <10 | 05/13/11 | 8270c & 625 | --- | 15.2 | 47.1 | 92.1 | 59 |
| Anthracene | <10 | µg/L | 10 | <10 | 05/13/11 | 8270c & 625 | --- | 16.1 | 58.8 | 93.4 | 82.7 |
| Benzidine | <40 | µg/L | 40 | <40 | 05/13/11 | 8270c & 625 | --- | 14.9 | 54.8 | 85.5 | 93.8 |
| Benzo[a]anthracene | <10 | µg/L | 10 | <10 | 05/13/11 | 8270c & 625 | --- | 11 | 56.4 | 93.7 | 84.6 |
| Benzo[a]pyrene | <10 | µg/L | 10 | <10 | 05/13/11 | 8270c & 625 | --- | 10.9 | 57.9 | 97 | 86 |
| Benzo[b]fluoranthene | <10 | µg/L | 10 | <10 | 05/13/11 | 8270c & 625 | --- | 13.2 | 55 | 97.6 | 81.8 |
| Benzo[g,h,i]perylene | <10 | µg/L | 10 | <10 | 05/13/11 | 8270c & 625 | --- | 12.2 | 64.6 | 90.3 | 101.1 |
| Benzo[j,k]fluoranthene | <10 | µg/L | 10 | <10 | 05/13/11 | 8270c & 625 | --- | 13.9 | 56.2 | 94.7 | 83.4 |
| Benzoic acid | <40 | µg/L | 40 | <40 | 05/13/11 | 8270c & 625 | --- | 26.7 | 1.9 | 88.3 | 4 |
| Benzyl alcohol | <20 | µg/L | 20 | <20 | 05/13/11 | 8270c & 625 | --- | 19.9 | 38 | 96.5 | 71.7 |
| bis(2-Chloroethoxy)methane | <10 | µg/L | 10 | <10 | 05/13/11 | 8270c & 625 | --- | 5 | 49.2 | 90.9 | 80.7 |
| bis(2-Chloroethyl)ether | <10 | µg/L | 10 | <10 | 05/13/11 | 8270c & 625 | --- | 6.5 | 44.7 | 93.2 | 66.3 |
| bis(2-chloroisopropyl)ether | <10 | µg/L | 10 | <10 | 05/13/11 | 8270c & 625 | --- | 9.8 | 40.2 | 90.9 | 61.4 |
| bis(2-Ethylhexyl)phthalate | <10 | µg/L | 10 | <10 | 05/13/11 | 8270c & 625 | --- | 14.1 | 53.2 | 98.3 | 83.1 |
| Butyl benzyl phthalate | <10 | µg/L | 10 | <10 | 05/13/11 | 8270c & 625 | --- | 6.9 | 50.9 | 95.3 | 83.8 |
| Chrysene | <10 | µg/L | 10 | <10 | 05/13/11 | 8270c & 625 | --- | 9.5 | 55.6 | 91.8 | 86.1 |
| Di-n-butyl phthalate (Dibutylphthalate) | <10 | µg/L | 10 | <10 | 05/13/11 | 8270c & 625 | --- | 15.5 | 54.3 | 91 | 80.4 |
| Di-n-octylphthalate (Dioctylphthalate) | <10 | µg/L | 10 | <10 | 05/13/11 | 8270c & 625 | --- | 11.9 | 52.7 | 101.8 | 82.1 |
| Dibenz[a,h]acridine | <10 | µg/L | 10 | <10 | 05/13/11 | 8270c & 625 | --- | 16.6 | 59.1 | 93.3 | 89.1 |
| Dibenz[a,h]anthracene | <10 | µg/L | 10 | <10 | 05/13/11 | 8270c & 625 | --- | 13.3 | 64.2 | 90.5 | 95.7 |
| Dibenzofuran | <10 | µg/L | 10 | <10 | 05/13/11 | 8270c & 625 | --- | 13 | 44.2 | 96 | 65.2 |
| Diethylphthalate | <10 | µg/L | 10 | <10 | 05/13/11 | 8270c & 625 | --- | 15 | 53.6 | 96.9 | 75.5 |
| Dimethylphthalate | <10 | µg/L | 10 | <10 | 05/13/11 | 8270c & 625 | --- | 9.2 | 53 | 94.9 | 77.6 |
| Fluoranthene | <10 | µg/L | 10 | <10 | 05/13/11 | 8270c & 625 | --- | 11.8 | 57.5 | 95.1 | 84.3 |
| Fluorene | <10 | µg/L | 10 | <10 | 05/13/11 | 8270c & 625 | J, | 9.2 | 49.9 | 97.1 | 72.3 |
| Hexachlorobenzene | <10 | µg/L | 10 | <10 | 05/13/11 | 8270c & 625 | --- | 12.4 | 55.9 | 93.6 | 75.6 |
| Hexachlorobutadiene | <10 | µg/L | 10 | <10 | 05/13/11 | 8270c & 625 | --- | 1.9 | 19.5 | 93 | 31.8 |

Client: Gainco, Inc.
Attn: Paul Supak

Project ID: Falcon Refinery
Sample Name: Tank 2

Report#/Lab ID#: 352493
Sample Matrix: water

REPORT OF ANALYSIS-cont.

QUALITY ASSURANCE DATA ¹

| Parameter | Result | Units | RQL ⁵ | Blank | Date | Method ⁶ | Data Qual. ⁷ | Prec. ² | Recov. ³ | CCV ⁴ | LCS ⁴ |
|-----------------------------------|--------|-------|------------------|-------|----------|---------------------|-------------------------|--------------------|---------------------|------------------|------------------|
| Hexachlorocyclopentadiene (HCCPD) | <10 | µg/L | 10 | <10 | 05/13/11 | 8270c & 625 | S1,M, | 0 | 0 | 102.8 | 14.8 |
| Hexachloroethane | <10 | µg/L | 10 | <10 | 05/13/11 | 8270c & 625 | --- | 2.9 | 19.7 | 94.8 | 31.4 |
| Indene | <10 | µg/L | 10 | <10 | 05/13/11 | 8270c & 625 | J,N, | 11.5 | 34.2 | 93.3 | 50.2 |
| Indeno[1,2,3-cd]pyrene | <10 | µg/L | 10 | <10 | 05/13/11 | 8270c & 625 | --- | 16.1 | 61 | 92.2 | 89.4 |
| Isophorone | <10 | µg/L | 10 | <10 | 05/13/11 | 8270c & 625 | --- | 6.7 | 57.6 | 90.4 | 95.3 |
| Methylchrysene | <10 | µg/L | 10 | <10 | 05/13/11 | 8270c & 625 | N, | 10.2 | 46.9 | 91.9 | 69.2 |
| N-Nitrosodi-n-propylamine | <10 | µg/L | 10 | <10 | 05/13/11 | 8270c & 625 | --- | 4.4 | 43.1 | 88.7 | 71.9 |
| N-Nitrosodimethylamine | <10 | µg/L | 10 | <10 | 05/13/11 | 8270c & 625 | --- | 5.9 | 31 | 94.3 | 41.5 |
| N-Nitrosodiphenylamine | <10 | µg/L | 10 | <10 | 05/13/11 | 8270c & 625 | --- | 14.4 | 59.5 | 91.7 | 85.4 |
| Naphthalene | 14.3 | µg/L | 10 | <10 | 05/13/11 | 8270c & 625 | --- | 5.5 | 31.7 | 90.7 | 49.8 |
| Nitrobenzene | <10 | µg/L | 10 | <10 | 05/13/11 | 8270c & 625 | --- | 1.6 | 44.1 | 91.2 | 72.3 |
| Pentachlorophenol | <10 | µg/L | 10 | <10 | 05/13/11 | 8270c & 625 | --- | 9.5 | 33.2 | 101.3 | 54.8 |
| Phenanthrene | <10 | µg/L | 10 | <10 | 05/13/11 | 8270c & 625 | --- | 12.2 | 59.8 | 92.6 | 86 |
| Phenol | <10 | µg/L | 10 | <10 | 05/13/11 | 8270c & 625 | J, | 15.4 | 17.9 | 93.1 | 25 |
| Pyrene | <10 | µg/L | 10 | <10 | 05/13/11 | 8270c & 625 | --- | 11.1 | 58.7 | 95.5 | 85.8 |
| Pyridine | <10 | µg/L | 10 | <10 | 05/13/11 | 8270c & 625 | --- | 19.5 | 26.1 | 94.6 | 34.1 |
| Quinoline | <10 | µg/L | 10 | <10 | 05/13/11 | 8270c & 625 | --- | 9.8 | 51.2 | 91.8 | 79.4 |

| | | |
|-----------------------------|------------------------------------|--------------------------------|
| Client: Gainco, Inc. | Project ID: Falcon Refinery | Report#/Lab ID#: 352493 |
| Attn: Paul Supak | Sample Name: Tank 2 | Sample Matrix: water |

REPORT OF SURROGATE RECOVERY

| Surrogate Compound | Method | Recovery | Recovery Limits | Date Analyzed | Data Qualifiers |
|-----------------------|-------------|----------|-----------------|---------------|-----------------|
| 1,2-Dichloroethane-d4 | 8260b & 624 | 85.6 | 70-125 | 05/12/11 | --- |
| 4-Bromofluorobenzene | 8260b & 624 | 96.7 | 80-115 | 05/12/11 | --- |
| Toluene-d8 | 8260b & 624 | 102.7 | 78-115 | 05/12/11 | --- |
| 2-Fluorobiphenyl | 8270c & 625 | 44.2 | 10-110 | 05/13/11 | --- |
| 2-Fluorophenol | 8270c & 625 | 26.6 | 10-110 | 05/13/11 | --- |
| 2,4,6-Tribromophenol | 8270c & 625 | 54 | 10-120 | 05/13/11 | --- |
| Nitrobenzene-d5 | 8270c & 625 | 42.4 | 10-110 | 05/13/11 | --- |
| Phenol-d6 | 8270c & 625 | 16.5 | 10-110 | 05/13/11 | --- |
| Terphenyl-d14 | 8270c & 625 | 27.7 | 10-115 | 05/13/11 | --- |

Data Qualifiers: D= Surrogates diluted and X= Surrogates outside advisory recovery limits.

Exceptions Report (FINAL SECTION / END-OF-REPORT):**Report #/Lab ID#:** 352493 **Matrix:** water**Client:** Gainco, Inc.**Attn:** Paul Supak**Project ID:** Falcon Refinery**Sample Name:** Tank 2

Unless otherwise identified by data qualifier "N" or by an exception report, all reported results represent parameters and tests for which AnalySys maintains NELAC certification; or results provided by a subcontractor with NELAC certification for the test results provided.

**Sample Temperature/Condition:** ≤6°C

The typical sample temperature criteria (except for metals by ICP, GFAA and AA and a very few other tests) is ≤ 6°C. Possible exceptions include samples submitted to laboratory within such a short time after sampling that cooling measures used in the field and during transport had insufficient time to achieve desired temperatures in the samples (see sample collection and sample receipt times) and samples where the temperature could not be measured due to sample submission in a manner precluding temperature measurement without impacting sample integrity (ex. in a bottle with no cooler).

Standard sample acceptability conditions met? : YES

Sample received in appropriate container(s), at appropriate temperature and pH.

J flag Discussion:

A J-flag data qualifier indicates that the raw calculated analyte concentration in the sample (uncorrected for background levels/blanks and other potential sources of sampling and analytical contamination), though less than the Reported Quantitation Limit (RQL) is greater than the Detection Limit. Because the reported result is below the quantitation limit for this project/sample (or test procedure), GC/MS organics results may or MAY NOT have been verified as to the presence and relative ratio of target ions (eg. the material causing the J flag "hit" in such situations may be nothing more than background ion-fragment noise.)

Comments pertaining to Data Qualifiers and QC data (where applicable):

| Parameter | Qualif. | Comments |
|--------------------------------------|---------|---|
| 1,1,1-Trichloroethane | L | Lab control sample (LCS or spiked blank). LCS recov-high (high bias). Sample result < MDL. No impact. |
| 1,1,1-Trichloroethane | C | Cont. Calib. Verification (CCV). CCV recov-high (high bias). Sample result < MDL. No impact. |
| 1,1,1-Trichloroethane | S | Spike (MS,MSD,PDS) recovery issue. MS, MSD & PDS recovery outside acceptance range. LCS fails or not available. Probable sample impact. |
| 1,2-Dichloropropane | C | Cont. Calib. Verification (CCV). CCV recov-high (high bias). Sample result < MDL. No impact. |
| Methylene chloride (Dichloromethane) | J | See J-flag discussion above. |
| 1-Methylnaphthalene | N | NELAC accreditation for this analyte not available from TCEQ. 30 TAC§25.6(4) applies. |
| 2-Methylnaphthalene | J | See J-flag discussion above. |
| 2-Methylphenol (o-Cresol) | J | See J-flag discussion above. |
| Fluorene | J | See J-flag discussion above. |
| Hexachlorocyclopentadiene (HCCPD) | S1 | Spike (MS,MSD) recovery issue. MS & MSD recovery outside acceptance range. LCS recovery OK. Probable matrix interference. |
| Indene | J | See J-flag discussion above. |
| Indene | N | NELAC accreditation for this analyte not available from TCEQ. 30 TAC§25.6(4) applies. |
| Methylchrysene | N | NELAC accreditation for this analyte not available from TCEQ. 30 TAC§25.6(4) applies. |
| Phenol | J | See J-flag discussion above. |

Client: Gainco, Inc.
Attn: Paul Supak
Address:

Phone: FAX:

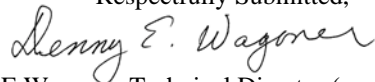
Report#/Lab ID#: 352494 Report Date: 05/16/11
Project ID: Falcon Refinery
Sample Name: Tank 27
Sample Matrix: water
Date Received: 05/10/2011 Time: 16:11
Date Sampled: 05/10/2011 Time: 12:00

REPORT OF ANALYSIS

QUALITY ASSURANCE DATA ¹

| Parameter | Result | Units | RQL ⁵ | Blank | Date | Method ⁶ | Data Qual. ⁷ | Prec. ² | Recov. ³ | CCV ⁴ | LCS ⁴ |
|--|---------|-------|------------------|---------|----------|---------------------|-------------------------|--------------------|---------------------|------------------|------------------|
| A/BN extraction-625/8270 | --- | --- | --- | --- | 05/13/11 | 3510 | --- | --- | --- | --- | --- |
| Metals Dig.-Hg | --- | --- | --- | --- | 05/11/11 | 7470 & SM3112B | --- | --- | --- | --- | --- |
| Metals Dig.-Total | --- | --- | --- | --- | 05/11/11 | 200.2 & 3005A | --- | --- | --- | --- | --- |
| Arsenic/ICPMS | <0.002 | mg/L | 0.002 | <0.002 | 05/13/11 | 6020A & 200.8 | J, | 0 | 99.5 | 98.3 | 99.6 |
| Barium/ICPMS | 0.209 | mg/L | 0.002 | <0.002 | 05/13/11 | 6020A & 200.8 | --- | 1 | 104.9 | 106.6 | 100 |
| Cadmium/ICPMS | <0.001 | mg/L | 0.001 | <0.001 | 05/13/11 | 6020A & 200.8 | --- | 0.8 | 93.9 | 100.5 | 92.7 |
| Chromium/ICPMS | <0.002 | mg/L | 0.002 | <0.002 | 05/13/11 | 6020A & 200.8 | --- | 1.5 | 99.5 | 98.5 | 100 |
| Lead/ICPMS | <0.002 | mg/L | 0.002 | <0.002 | 05/13/11 | 6020A & 200.8 | J, | 0.1 | 87 | 95.6 | 92.2 |
| Mercury/CVAA | <0.0002 | mg/L | 0.0002 | <0.0002 | 05/12/11 | 7470&SM3112B | --- | 3.08 | 90 | 107 | 107 |
| Selenium/ICPMS | <0.002 | mg/L | 0.002 | <0.002 | 05/13/11 | 6020A & 200.8 | --- | 1.7 | 106.2 | 104 | 106.9 |
| Silver/ICPMS | <0.001 | mg/L | 0.001 | <0.001 | 05/13/11 | 6020A & 200.8 | --- | 1 | 92.7 | 97.6 | 92.2 |
| Volatile organics-8260 | --- | --- | --- | --- | 05/12/11 | 8260b(5030/5035) | --- | --- | --- | --- | --- |
| 1,1-Dichloroethane | <1 | µg/L | 1 | <1 | 05/12/11 | 8260b & 624 | --- | 5.9 | 136.5 | 119.2 | 121.3 |
| 1,1-Dichloroethene | <1 | µg/L | 1 | <1 | 05/12/11 | 8260b & 624 | --- | 7.3 | 107.9 | 94.5 | 100.3 |
| 1,1,1-Trichloroethane | <1 | µg/L | 1 | <1 | 05/12/11 | 8260b & 624 | C,L,S, | 8.2 | 143.9 | 127.9 | 125.8 |
| 1,1,1,2-Tetrachloroethane | <1 | µg/L | 1 | <1 | 05/12/11 | 8260b & 624 | --- | 1 | 107.3 | 101.8 | 102.7 |
| 1,1,2-Trichloroethane | <1 | µg/L | 1 | <1 | 05/12/11 | 8260b & 624 | --- | 2.2 | 124.7 | 116.1 | 114.5 |
| 1,1,2,2-Tetrachloroethane | <1 | µg/L | 1 | <1 | 05/12/11 | 8260b & 624 | --- | 1.6 | 129.1 | 128.8 | 122.8 |
| 1,2-Dibromo-3-chloropropane | <2 | µg/L | 2 | <2 | 05/12/11 | 8260b & 624 | --- | 2.6 | 112.3 | 104.9 | 106.1 |
| 1,2-Dibromoethane (Ethylene dibromide) | <1 | µg/L | 1 | <1 | 05/12/11 | 8260b & 624 | --- | 0.5 | 118.1 | 112 | 112.5 |
| 1,2-Dichlorobenzene | <1 | µg/L | 1 | <1 | 05/12/11 | 8260b & 624 | --- | 1.2 | 111.4 | 110.8 | 110 |

This analytical report is respectfully submitted by AnalySys, Inc. The enclosed results reflect only the sample identified above. The results have been carefully reviewed and to the best of my knowledge, unless otherwise indicated, meet NELAC requirements as described by AnalySys, Inc.'s Quality Assurance/Quality Control Program. © Copyright 2003, AnalySys, Inc., Austin, TX. All rights reserved. No part of this publication may be reproduced or transmitted in any form or by any means without the express written consent of AnalySys, Inc.

Respectfully Submitted,

D.E. Wagoner, Technical Director (or designee)

1. Quality assurance data for the sample batch which included this sample. 2. Precision (PREC) is the absolute value of the relative percent difference between duplicate results. 3. Recovery (Recov.) is the percent of analyte recovered from a spiked sample. 4. Calibration Verification (CCV) and Laboratory Control Sample (LCS) results are expressed as the percent recovery of analyte. 5. Reporting Quantitation Limits (RQL), typically at or above the Practical Quantitation Limit (PQL) of the analytical method. 6. Method numbers typically denote USEPA procedures. Less than ("<") values reflect nominal quantitation limits adjusted for any required dilutions. 7. Data Qualifiers are J = analyte detected between the RQL and the MDL. B = Analyte detected in associated method blank (s). C=poor CCV recovery. L=poor LCS recovery. S & S1 =MS and/or MSD recovery exceed advisory limits. S2 =Post digestion spike (PDS) recovery exceeds advisory limit. S3 =MS and/or MSD and PDS recoveries exceed advisory limits. P =Precision higher than advisory limit. M =Matrix interference. N=not NELACcertified. N1=subcontract result enquire concerning NELAC certification. Solid sample results for all metals, except Mercury, reported on a dry weight basis (DWB)s. All other results for solid samples reported on an as received basis unless specifically identified as DWB.

Client: Gainco, Inc.
Attn: Paul Supak

Project ID: Falcon Refinery
Sample Name: Tank 27

Report#/Lab ID#: 352494
Sample Matrix: water

REPORT OF ANALYSIS-cont.
QUALITY ASSURANCE DATA ¹

| Parameter | Result | Units | RQL ⁵ | Blank | Date | Method ⁶ | Data Qual. ⁷ | Prec. ² | Recov. ³ | CCV ⁴ | LCS ⁴ |
|--------------------------------------|--------|-------|------------------|-------|----------|---------------------|-------------------------|--------------------|---------------------|------------------|------------------|
| 1,2-Dichloroethane | <1 | µg/L | 1 | <1 | 05/12/11 | 8260b & 624 | --- | 0.8 | 103.3 | 93.3 | 90.9 |
| 1,2-Dichloropropane | <1 | µg/L | 1 | <1 | 05/12/11 | 8260b & 624 | C, | 2 | 132.1 | 120.6 | 117.1 |
| 1,3-Dichlorobenzene | <1 | µg/L | 1 | <1 | 05/12/11 | 8260b & 624 | --- | 0.1 | 109.7 | 108.2 | 107.9 |
| 1,4-Dichlorobenzene | <1 | µg/L | 1 | <1 | 05/12/11 | 8260b & 624 | --- | 1 | 108.5 | 106.9 | 105.3 |
| 1,4-Dioxane | <20 | µg/L | 20 | <20 | 05/12/11 | 8260b & 624 | --- | 2.5 | 119.4 | 105 | 118.4 |
| 2-Butanone (MEK) | <2 | µg/L | 2 | <2 | 05/12/11 | 8260b & 624 | --- | 2.6 | 137.8 | 115.3 | 128.2 |
| 2-Chloroethyl vinyl ether | <1 | µg/L | 1 | <1 | 05/12/11 | 8260b & 624 | --- | 16.2 | 108.9 | 93.4 | 100.4 |
| 2-Hexanone | <2 | µg/L | 2 | <2 | 05/12/11 | 8260b & 624 | --- | 0.6 | 115.7 | 111.4 | 111.7 |
| 4-Methyl-2-pentanone (MIBK) | <2 | µg/L | 2 | <2 | 05/12/11 | 8260b & 624 | --- | 0.8 | 129.4 | 114.8 | 119.6 |
| Acetone (2-Propanone) | <10 | µg/L | 10 | <10 | 05/12/11 | 8260b & 624 | J, | 5.5 | 105.3 | 92.3 | 96 |
| Acetonitrile | <20 | µg/L | 20 | <20 | 05/12/11 | 8260b & 624 | --- | 1.2 | 133.3 | 112.4 | 118.1 |
| Acrolein | <10 | µg/L | 10 | <10 | 05/12/11 | 8260b & 624 | --- | 3.9 | 123 | 106.5 | 107.7 |
| Acrylonitrile | <10 | µg/L | 10 | <10 | 05/12/11 | 8260b & 624 | --- | 0.4 | 139.8 | 122.9 | 123.3 |
| Benzene | <1 | µg/L | 1 | <1 | 05/12/11 | 8260b & 624 | J, | 4.5 | 126.1 | 114.9 | 111.1 |
| Bromobenzene | <1 | µg/L | 1 | <1 | 05/12/11 | 8260b & 624 | --- | 1 | 117.2 | 116 | 111.7 |
| Bromodichloromethane | <1 | µg/L | 1 | <1 | 05/12/11 | 8260b & 624 | --- | 2 | 110.1 | 99.8 | 98.2 |
| Bromoform (Tribromomethane) | <1 | µg/L | 1 | <1 | 05/12/11 | 8260b & 624 | --- | 0.9 | 100.9 | 95.4 | 95.4 |
| Bromomethane (Methyl bromide) | <2 | µg/L | 2 | <2 | 05/12/11 | 8260b & 624 | --- | 4.7 | 91.6 | 80.7 | 86.8 |
| Carbon disulfide | <2 | µg/L | 2 | <2 | 05/12/11 | 8260b & 624 | --- | 7.9 | 127.6 | 113.6 | 128.7 |
| Carbon tetrachloride | <1 | µg/L | 1 | <1 | 05/12/11 | 8260b & 624 | --- | 5 | 97.6 | 89.7 | 89.7 |
| Chlorobenzene | <1 | µg/L | 1 | <1 | 05/12/11 | 8260b & 624 | --- | 1 | 113.7 | 107.7 | 108.4 |
| Chloroethane | <2 | µg/L | 2 | <2 | 05/12/11 | 8260b & 624 | --- | 8.1 | 106.7 | 96.2 | 92.8 |
| Chloroform (Trichloromethane) | <1 | µg/L | 1 | <1 | 05/12/11 | 8260b & 624 | --- | 2.2 | 120.9 | 107.8 | 110 |
| Chloromethane (Methyl chloride) | <2 | µg/L | 2 | <2 | 05/12/11 | 8260b & 624 | --- | 7.3 | 102.7 | 92 | 94.3 |
| cis-1,2-Dichloroethene | <1 | µg/L | 1 | <1 | 05/12/11 | 8260b & 624 | --- | 2.6 | 130.4 | 116.9 | 113.9 |
| cis-1,3-Dichloropropene | <1 | µg/L | 1 | <1 | 05/12/11 | 8260b & 624 | --- | 0.8 | 117.6 | 108 | 104.8 |
| Dibromochloromethane | <1 | µg/L | 1 | <1 | 05/12/11 | 8260b & 624 | --- | 2.7 | 107.5 | 104.3 | 103.9 |
| Dibromomethane (Methylene bromide) | <1 | µg/L | 1 | <1 | 05/12/11 | 8260b & 624 | --- | 2.3 | 116.5 | 103.4 | 101.6 |
| Dichlorodifluoromethane | <1 | µg/L | 1 | <1 | 05/12/11 | 8260b & 624 | --- | 11.8 | 52.2 | 48.2 | 45.7 |
| Ethylbenzene | <1 | µg/L | 1 | <1 | 05/12/11 | 8260b & 624 | --- | 1 | 114.6 | 109.3 | 109.2 |
| Iodomethane (Methyl iodide) | <2 | µg/L | 2 | <2 | 05/12/11 | 8260b & 624 | --- | 4.8 | 86.2 | 76 | 84.2 |
| m,p-Xylenes | <2 | µg/L | 2 | <2 | 05/12/11 | 8260b & 624 | --- | 0.9 | 113 | 107.6 | 107.6 |
| Methylene chloride (Dichloromethane) | <2 | µg/L | 2 | <2 | 05/12/11 | 8260b & 624 | J, | 2.5 | 137.5 | 113.8 | 116.4 |
| MTBE | 8.43 | µg/L | 1 | <1 | 05/12/11 | 8260b & 624 | --- | 4.2 | 135.2 | 118.3 | 113.3 |

Client: Gainco, Inc.
Attn: Paul Supak

Project ID: Falcon Refinery
Sample Name: Tank 27

Report#/Lab ID#: 352494
Sample Matrix: water

REPORT OF ANALYSIS-cont.
QUALITY ASSURANCE DATA ¹

| Parameter | Result | Units | RQL ⁵ | Blank | Date | Method ⁶ | Data Qual. ⁷ | Prec. ² | Recov. ³ | CCV ⁴ | LCS ⁴ |
|---------------------------------------|--------|-------|------------------|-------|----------|---------------------|-------------------------|--------------------|---------------------|------------------|------------------|
| o-Xylene | <1 | µg/L | 1 | <1 | 05/12/11 | 8260b & 624 | --- | 0.7 | 114.6 | 107.5 | 109.5 |
| Styrene | <1 | µg/L | 1 | <1 | 05/12/11 | 8260b & 624 | --- | 1.2 | 116.1 | 111 | 110.4 |
| Tetrachloroethene (Perchloroethylene) | <1 | µg/L | 1 | <1 | 05/12/11 | 8260b & 624 | --- | 4.1 | 105.6 | 99.4 | 100.1 |
| Toluene | <1 | µg/L | 1 | <1 | 05/12/11 | 8260b & 624 | --- | 3.7 | 120.5 | 108.7 | 107.4 |
| trans-1,2-Dichloroethene | <1 | µg/L | 1 | <1 | 05/12/11 | 8260b & 624 | --- | 5.3 | 127 | 113.8 | 123.9 |
| trans-1,3-Dichloropropene | <1 | µg/L | 1 | <1 | 05/12/11 | 8260b & 624 | --- | 1.7 | 112.2 | 108.7 | 112.2 |
| Trichloroethene | <1 | µg/L | 1 | <1 | 05/12/11 | 8260b & 624 | --- | 3.1 | 111.7 | 104 | 100.4 |
| Trichlorofluoromethane | <1 | µg/L | 1 | <1 | 05/12/11 | 8260b & 624 | --- | 7.1 | 99.7 | 87.2 | 93.3 |
| Vinyl acetate | <2 | µg/L | 2 | <2 | 05/12/11 | 8260b & 624 | --- | 4.3 | 140.5 | 119.7 | 120.1 |
| Vinyl chloride | <1 | µg/L | 1 | <1 | 05/12/11 | 8260b & 624 | --- | 11 | 95.8 | 85.9 | 84.6 |
| Extractable organics-625/8270 | --- | --- | --- | --- | 05/13/11 | 8270c & 625 | --- | --- | --- | --- | --- |
| 1-Methylnaphthalene | <10 | µg/L | 10 | <10 | 05/13/11 | 8270c & 625 | N, | 8 | 28.8 | 88.1 | 45.2 |
| 1,2-Diphenylhydrazine | <10 | µg/L | 10 | <10 | 05/13/11 | 8270c & 625 | --- | 15.7 | 55.8 | 92.5 | 78.6 |
| 1,2,4-Trichlorobenzene | <10 | µg/L | 10 | <10 | 05/13/11 | 8270c & 625 | --- | 3.3 | 21.6 | 90 | 37.6 |
| 2-Chloronaphthalene | <10 | µg/L | 10 | <10 | 05/13/11 | 8270c & 625 | --- | 9 | 29.1 | 95.2 | 46.2 |
| 2-Chlorophenol | <10 | µg/L | 10 | <10 | 05/13/11 | 8270c & 625 | --- | 7.8 | 40.2 | 93.2 | 57.1 |
| 2-Methylnaphthalene | <10 | µg/L | 10 | <10 | 05/13/11 | 8270c & 625 | --- | 9.8 | 27.9 | 90.4 | 43.3 |
| 2-Methylphenol (o-Cresol) | <10 | µg/L | 10 | <10 | 05/13/11 | 8270c & 625 | --- | 22.2 | 46.5 | 93.9 | 63.6 |
| 2-Nitroaniline | <50 | µg/L | 50 | <50 | 05/13/11 | 8270c & 625 | --- | 9.2 | 52.6 | 93.9 | 81.5 |
| 2-Nitrophenol | <10 | µg/L | 10 | <10 | 05/13/11 | 8270c & 625 | --- | 5.9 | 36.6 | 98 | 56.1 |
| 2,4-Dichlorophenol | <10 | µg/L | 10 | <10 | 05/13/11 | 8270c & 625 | --- | 6.5 | 41.4 | 95.9 | 65.9 |
| 2,4-Dimethylphenol | <10 | µg/L | 10 | <10 | 05/13/11 | 8270c & 625 | --- | 3.8 | 48.3 | 86.8 | 78.3 |
| 2,4-Dinitrophenol | <10 | µg/L | 10 | <10 | 05/13/11 | 8270c & 625 | --- | 18.1 | 19.8 | 98.7 | 34.6 |
| 2,4-Dinitrotoluene | <10 | µg/L | 10 | <10 | 05/13/11 | 8270c & 625 | --- | 13.3 | 52.2 | 102.9 | 77.7 |
| 2,4,6-Trichlorophenol | <10 | µg/L | 10 | <10 | 05/13/11 | 8270c & 625 | --- | 10.4 | 39.4 | 96.4 | 56 |
| 2,6-Dichlorophenol | <10 | µg/L | 10 | <10 | 05/13/11 | 8270c & 625 | --- | 3.3 | 36.6 | 92.3 | 52.6 |
| 2,6-Dinitrotoluene | <10 | µg/L | 10 | <10 | 05/13/11 | 8270c & 625 | --- | 16.2 | 55.8 | 97.8 | 82.3 |
| 3-Nitroaniline | <50 | µg/L | 50 | <50 | 05/13/11 | 8270c & 625 | --- | 8.7 | 50.9 | 102 | 82.9 |
| 3,3'-Dichlorobenzidine | <20 | µg/L | 20 | <20 | 05/13/11 | 8270c & 625 | --- | 13.3 | 56.2 | 96.6 | 86.9 |
| 3&4 Methylphenol (m&p-Cresol) | <20 | µg/L | 20 | <20 | 05/13/11 | 8270c & 625 | --- | 15.8 | 40.8 | 96 | 59.6 |
| 4-Bromophenyl phenyl ether | <10 | µg/L | 10 | <10 | 05/13/11 | 8270c & 625 | --- | 19.5 | 52.4 | 95.6 | 70.8 |
| 4-Chloro-3-methylphenol | <20 | µg/L | 20 | <20 | 05/13/11 | 8270c & 625 | --- | 4.1 | 51.6 | 92.8 | 73.9 |
| 4-Chloroaniline (p-Chloroaniline) | <20 | µg/L | 20 | <20 | 05/13/11 | 8270c & 625 | --- | 15.7 | 36.3 | 106.7 | 77.6 |

Client: Gainco, Inc.
Attn: Paul Supak

Project ID: Falcon Refinery
Sample Name: Tank 27

Report#/Lab ID#: 352494
Sample Matrix: water

REPORT OF ANALYSIS-cont.
QUALITY ASSURANCE DATA ¹

| Parameter | Result | Units | RQL ⁵ | Blank | Date | Method ⁶ | Data Qual. ⁷ | Prec. ² | Recov. ³ | CCV ⁴ | LCS ⁴ |
|---|--------|-------|------------------|-------|----------|---------------------|-------------------------|--------------------|---------------------|------------------|------------------|
| 4-Chlorophenyl phenyl ether | <10 | µg/L | 10 | <10 | 05/13/11 | 8270c & 625 | --- | 13.2 | 45.8 | 96.5 | 66.6 |
| 4-Nitroaniline | <20 | µg/L | 20 | <20 | 05/13/11 | 8270c & 625 | --- | 6.7 | 44.8 | 97.5 | 73.3 |
| 4-Nitrophenol | <50 | µg/L | 50 | <50 | 05/13/11 | 8270c & 625 | J, | 7.2 | 12.6 | 83 | 16.2 |
| 4,6-Dinitro-2-methylphenol | <25 | µg/L | 25 | <25 | 05/13/11 | 8270c & 625 | --- | 19.5 | 35.5 | 101.9 | 50.3 |
| 7,12-Dimethylbenz[a]anthracene | <10 | µg/L | 10 | <10 | 05/13/11 | 8270c & 625 | --- | 14.5 | 56.9 | 99.4 | 84.4 |
| Acenaphthene | <10 | µg/L | 10 | <10 | 05/13/11 | 8270c & 625 | --- | 12.8 | 40.6 | 94.3 | 58.1 |
| Acenaphthylene | <10 | µg/L | 10 | <10 | 05/13/11 | 8270c & 625 | --- | 8.4 | 40.9 | 96.1 | 61.8 |
| Aniline | <10 | µg/L | 10 | <10 | 05/13/11 | 8270c & 625 | --- | 15.2 | 47.1 | 92.1 | 59 |
| Anthracene | <10 | µg/L | 10 | <10 | 05/13/11 | 8270c & 625 | --- | 16.1 | 58.8 | 93.4 | 82.7 |
| Benzidine | <40 | µg/L | 40 | <40 | 05/13/11 | 8270c & 625 | --- | 14.9 | 54.8 | 85.5 | 93.8 |
| Benzo[a]anthracene | <10 | µg/L | 10 | <10 | 05/13/11 | 8270c & 625 | --- | 11 | 56.4 | 93.7 | 84.6 |
| Benzo[a]pyrene | <10 | µg/L | 10 | <10 | 05/13/11 | 8270c & 625 | --- | 10.9 | 57.9 | 97 | 86 |
| Benzo[b]fluoranthene | <10 | µg/L | 10 | <10 | 05/13/11 | 8270c & 625 | --- | 13.2 | 55 | 97.6 | 81.8 |
| Benzo[g,h,i]perylene | <10 | µg/L | 10 | <10 | 05/13/11 | 8270c & 625 | --- | 12.2 | 64.6 | 90.3 | 101.1 |
| Benzo[j,k]fluoranthene | <10 | µg/L | 10 | <10 | 05/13/11 | 8270c & 625 | --- | 13.9 | 56.2 | 94.7 | 83.4 |
| Benzoic acid | <40 | µg/L | 40 | <40 | 05/13/11 | 8270c & 625 | --- | 26.7 | 1.9 | 88.3 | 4 |
| Benzyl alcohol | <20 | µg/L | 20 | <20 | 05/13/11 | 8270c & 625 | --- | 19.9 | 38 | 96.5 | 71.7 |
| bis(2-Chloroethoxy)methane | <10 | µg/L | 10 | <10 | 05/13/11 | 8270c & 625 | --- | 5 | 49.2 | 90.9 | 80.7 |
| bis(2-Chloroethyl)ether | <10 | µg/L | 10 | <10 | 05/13/11 | 8270c & 625 | --- | 6.5 | 44.7 | 93.2 | 66.3 |
| bis(2-chloroisopropyl)ether | <10 | µg/L | 10 | <10 | 05/13/11 | 8270c & 625 | --- | 9.8 | 40.2 | 90.9 | 61.4 |
| bis(2-Ethylhexyl)phthalate | <10 | µg/L | 10 | <10 | 05/13/11 | 8270c & 625 | --- | 14.1 | 53.2 | 98.3 | 83.1 |
| Butyl benzyl phthalate | <10 | µg/L | 10 | <10 | 05/13/11 | 8270c & 625 | --- | 6.9 | 50.9 | 95.3 | 83.8 |
| Chrysene | <10 | µg/L | 10 | <10 | 05/13/11 | 8270c & 625 | --- | 9.5 | 55.6 | 91.8 | 86.1 |
| Di-n-butyl phthalate (Dibutylphthalate) | <10 | µg/L | 10 | <10 | 05/13/11 | 8270c & 625 | --- | 15.5 | 54.3 | 91 | 80.4 |
| Di-n-octylphthalate (Dioctylphthalate) | <10 | µg/L | 10 | <10 | 05/13/11 | 8270c & 625 | --- | 11.9 | 52.7 | 101.8 | 82.1 |
| Dibenz[a,h]acridine | <10 | µg/L | 10 | <10 | 05/13/11 | 8270c & 625 | --- | 16.6 | 59.1 | 93.3 | 89.1 |
| Dibenz[a,h]anthracene | <10 | µg/L | 10 | <10 | 05/13/11 | 8270c & 625 | --- | 13.3 | 64.2 | 90.5 | 95.7 |
| Dibenzofuran | <10 | µg/L | 10 | <10 | 05/13/11 | 8270c & 625 | --- | 13 | 44.2 | 96 | 65.2 |
| Diethylphthalate | <10 | µg/L | 10 | <10 | 05/13/11 | 8270c & 625 | --- | 15 | 53.6 | 96.9 | 75.5 |
| Dimethylphthalate | <10 | µg/L | 10 | <10 | 05/13/11 | 8270c & 625 | --- | 9.2 | 53 | 94.9 | 77.6 |
| Fluoranthene | <10 | µg/L | 10 | <10 | 05/13/11 | 8270c & 625 | --- | 11.8 | 57.5 | 95.1 | 84.3 |
| Fluorene | <10 | µg/L | 10 | <10 | 05/13/11 | 8270c & 625 | --- | 9.2 | 49.9 | 97.1 | 72.3 |
| Hexachlorobenzene | <10 | µg/L | 10 | <10 | 05/13/11 | 8270c & 625 | --- | 12.4 | 55.9 | 93.6 | 75.6 |
| Hexachlorobutadiene | <10 | µg/L | 10 | <10 | 05/13/11 | 8270c & 625 | --- | 1.9 | 19.5 | 93 | 31.8 |

Client: Gainco, Inc.
Attn: Paul Supak

Project ID: Falcon Refinery
Sample Name: Tank 27

Report#/Lab ID#: 352494
Sample Matrix: water

REPORT OF ANALYSIS-cont.
QUALITY ASSURANCE DATA ¹

| Parameter | Result | Units | RQL ⁵ | Blank | Date | Method ⁶ | Data Qual. ⁷ | Prec. ² | Recov. ³ | CCV ⁴ | LCS ⁴ |
|-----------------------------------|--------|-------|------------------|-------|----------|---------------------|-------------------------|--------------------|---------------------|------------------|------------------|
| Hexachlorocyclopentadiene (HCCPD) | <10 | µg/L | 10 | <10 | 05/13/11 | 8270c & 625 | S1,M, | 0 | 0 | 102.8 | 14.8 |
| Hexachloroethane | <10 | µg/L | 10 | <10 | 05/13/11 | 8270c & 625 | --- | 2.9 | 19.7 | 94.8 | 31.4 |
| Indene | <10 | µg/L | 10 | <10 | 05/13/11 | 8270c & 625 | N, | 11.5 | 34.2 | 93.3 | 50.2 |
| Indeno[1,2,3-cd]pyrene | <10 | µg/L | 10 | <10 | 05/13/11 | 8270c & 625 | --- | 16.1 | 61 | 92.2 | 89.4 |
| Isophorone | <10 | µg/L | 10 | <10 | 05/13/11 | 8270c & 625 | --- | 6.7 | 57.6 | 90.4 | 95.3 |
| Methylchrysene | <10 | µg/L | 10 | <10 | 05/13/11 | 8270c & 625 | N, | 10.2 | 46.9 | 91.9 | 69.2 |
| N-Nitrosodi-n-propylamine | <10 | µg/L | 10 | <10 | 05/13/11 | 8270c & 625 | --- | 4.4 | 43.1 | 88.7 | 71.9 |
| N-Nitrosodimethylamine | <10 | µg/L | 10 | <10 | 05/13/11 | 8270c & 625 | --- | 5.9 | 31 | 94.3 | 41.5 |
| N-Nitrosodiphenylamine | <10 | µg/L | 10 | <10 | 05/13/11 | 8270c & 625 | --- | 14.4 | 59.5 | 91.7 | 85.4 |
| Naphthalene | <10 | µg/L | 10 | <10 | 05/13/11 | 8270c & 625 | --- | 5.5 | 31.7 | 90.7 | 49.8 |
| Nitrobenzene | <10 | µg/L | 10 | <10 | 05/13/11 | 8270c & 625 | --- | 1.6 | 44.1 | 91.2 | 72.3 |
| Pentachlorophenol | <10 | µg/L | 10 | <10 | 05/13/11 | 8270c & 625 | --- | 9.5 | 33.2 | 101.3 | 54.8 |
| Phenanthrene | <10 | µg/L | 10 | <10 | 05/13/11 | 8270c & 625 | --- | 12.2 | 59.8 | 92.6 | 86 |
| Phenol | <10 | µg/L | 10 | <10 | 05/13/11 | 8270c & 625 | --- | 15.4 | 17.9 | 93.1 | 25 |
| Pyrene | <10 | µg/L | 10 | <10 | 05/13/11 | 8270c & 625 | --- | 11.1 | 58.7 | 95.5 | 85.8 |
| Pyridine | <10 | µg/L | 10 | <10 | 05/13/11 | 8270c & 625 | --- | 19.5 | 26.1 | 94.6 | 34.1 |
| Quinoline | <10 | µg/L | 10 | <10 | 05/13/11 | 8270c & 625 | --- | 9.8 | 51.2 | 91.8 | 79.4 |

| | | |
|-----------------------------|------------------------------------|--------------------------------|
| Client: Gainco, Inc. | Project ID: Falcon Refinery | Report#/Lab ID#: 352494 |
| Attn: Paul Supak | Sample Name: Tank 27 | Sample Matrix: water |

REPORT OF SURROGATE RECOVERY

| Surrogate Compound | Method | Recovery | Recovery Limits | Date Analyzed | Data Qualifiers |
|-----------------------|-------------|----------|-----------------|---------------|-----------------|
| 1,2-Dichloroethane-d4 | 8260b & 624 | 82.8 | 70-125 | 05/12/11 | --- |
| 4-Bromofluorobenzene | 8260b & 624 | 97.8 | 80-115 | 05/12/11 | --- |
| Toluene-d8 | 8260b & 624 | 104.1 | 78-115 | 05/12/11 | --- |
| 2-Fluorobiphenyl | 8270c & 625 | 34.5 | 10-110 | 05/13/11 | --- |
| 2-Fluorophenol | 8270c & 625 | 24 | 10-110 | 05/13/11 | --- |
| 2,4,6-Tribromophenol | 8270c & 625 | 50.3 | 10-120 | 05/13/11 | --- |
| Nitrobenzene-d5 | 8270c & 625 | 37.3 | 10-110 | 05/13/11 | --- |
| Phenol-d6 | 8270c & 625 | 14.8 | 10-110 | 05/13/11 | --- |
| Terphenyl-d14 | 8270c & 625 | 43.2 | 10-115 | 05/13/11 | --- |

Data Qualifiers: D= Surrogates diluted and X= Surrogates outside advisory recovery limits.

Exceptions Report (FINAL SECTION / END-OF-REPORT):**Report #/Lab ID#:** 352494 **Matrix:** water**Client:** Gainco, Inc.**Attn:** Paul Supak**Project ID:** Falcon Refinery**Sample Name:** Tank 27

Unless otherwise identified by data qualifier "N" or by an exception report, all reported results represent parameters and tests for which AnalySys maintains NELAC certification; or results provided by a subcontractor with NELAC certification for the test results provided.

**Sample Temperature/Condition:** ≤6°C

The typical sample temperature criteria (except for metals by ICP, GFAA and AA and a very few other tests) is ≤ 6°C. Possible exceptions include samples submitted to laboratory within such a short time after sampling that cooling measures used in the field and during transport had insufficient time to achieve desired temperatures in the samples (see sample collection and sample receipt times) and samples where the temperature could not be measured due to sample submission in a manner precluding temperature measurement without impacting sample integrity (ex. in a bottle with no cooler).

Standard sample acceptability conditions met? : YES

Sample received in appropriate container(s), at appropriate temperature and pH.

J flag Discussion:

A J-flag data qualifier indicates that the raw calculated analyte concentration in the sample (uncorrected for background levels/blanks and other potential sources of sampling and analytical contamination), though less than the Reported Quantitation Limit (RQL) is greater than the Detection Limit. Because the reported result is below the quantitation limit for this project/sample (or test procedure), GC/MS organics results may or MAY NOT have been verified as to the presence and relative ratio of target ions (eg. the material causing the J flag "hit" in such situations may be nothing more than background ion-fragment noise.)

Comments pertaining to Data Qualifiers and QC data (where applicable):

| Parameter | Qualif. | Comments |
|--------------------------------------|---------|---|
| Arsenic/ICPMS | J | See J-flag discussion above. |
| Lead/ICPMS | J | See J-flag discussion above. |
| 1,1,1-Trichloroethane | L | Lab control sample (LCS or spiked blank). LCS recov-high (high bias). Sample result < MDL. No impact. |
| 1,1,1-Trichloroethane | C | Cont. Calib. Verification (CCV). CCV recov-high (high bias). Sample result < MDL. No impact. |
| 1,1,1-Trichloroethane | S | Spike (MS,MSD,PDS) recovery issue. MS, MSD & PDS recovery outside acceptance range. LCS fails or not available. Probable sample impact. |
| 1,2-Dichloropropane | C | Cont. Calib. Verification (CCV). CCV recov-high (high bias). Sample result < MDL. No impact. |
| Acetone (2-Propanone) | J | See J-flag discussion above. |
| Benzene | J | See J-flag discussion above. |
| Methylene chloride (Dichloromethane) | J | See J-flag discussion above. |
| 1-Methylnaphthalene | N | NELAC accreditation for this analyte not available from TCEQ. 30 TAC§25.6(4) applies. |
| 4-Nitrophenol | J | See J-flag discussion above. |
| Hexachlorocyclopentadiene (HCCPD) | S1 | Spike (MS,MSD) recovery issue. MS & MSD recovery outside acceptance range. LCS recovery OK. Probable matrix interference. |
| Indene | N | NELAC accreditation for this analyte not available from TCEQ. 30 TAC§25.6(4) applies. |
| Methylchrysene | N | NELAC accreditation for this analyte not available from TCEQ. 30 TAC§25.6(4) applies. |

Send Reports To:

ANALYSIS
INC.

2209 N. Padre Island Drive, Suite K
Corpus Christi, TX 78408
Ph (361) 289-6384 • Fax (361) 289-0875

Bill To (if different):

Company Name SAME
Address _____
City _____ State _____ Zip _____
ATTN: _____
Phone _____ Fax _____

No. of Containers and
Preservative
(TRRP-13 Mandatory)

Matrix

Analyze For

21

[illegible]

| |
|--|
| Special Instructions (such as special QC requirements, lists, methods, etc...) |
|--|

Email results to Paul Supak at psupak@gaincoinc.com

24 HR TAT

Temperature
upon receipt
(Consistent with
NELAC sec.
5.11) ($>0-6^{\circ}\text{C}$)

1.70C

(1) Unless specifically requested otherwise on this Chain-of-custody and/or attached documentation, all analyses will be conducted using ASI's method of choice and all data will be reported to ASI's normal reporting limits (MDL/PQL). For GC/MS volatiles and extractables, unless specific analytical parameter lists are specified on this chain-of-custody or attached to this chain-of-custody, ASI will default to Priority Pollutants or ASI's HSL list at ASI's option. Specific compound lists must be supplied for all GC procedures.

upon receipt
(Consistent with
NELAC sec.
5.11) (>0-6°C)

| Sample Relinquished By | | | | Sample Received By | | | | S.M. (2000) | |
|------------------------|-------------|---------|-------|--------------------|-------------|---------|-------|-------------|---|
| Name | Affiliation | Date | Time | Name | Affiliation | Date | Time | YES | X |
| Paul Spitz | GAINCO | 5/10/11 | 16:11 | C. Adams | ASI | 5/10/11 | 16:11 | NO | |

[Tendering of above described samples to AnalySys, Inc. for analytical testing constitutes agreement by buyer/sampler to AnalySys, Inc.'s standard terms.]

PRI PPD

10MAY11 05:45P

** LABEL **

Schd: VLP 0834

GLI 3061984112

Pcs: 3 of 5

AUSTIN, TX



From: ANALYSYS INC AUSTIN TX
512-385-5886

RECV: ANALYSYS

3512 MONTOPOLIS DR

2.4

Manual Wght:

227.5

Tariff Wght:

230.0

AUSTIN, TX 78751

Phone: 222-222-2222

PO/Ref #:

Priority

Agency Phone: (512) 454-9686

WWW.SHIPGREYHOUND.COM

☐ RUSH ¹³5-12-11 Date Due
₂₃

ASI-0012

Rev. 2

Prepared: 07/09/2009

Effective: 07/20/2009

ASI Sample Evaluation and Comment Tracking

Sample #'s: 352492-494 Client: Grainco Date: 5/10/11

ASI Proj.#: 3269S Proj. Name: Falcon Refinery # of C-O-C's: 1

In compliance with the NELAC standard, ASI is notifying you that the **SAMPLES** identified here and on the attached Chain-of-Custody were received by AnalySys, Inc. (ASI) with the following **INTEGRITY ISSUES** (any NO responses indicated below). In order to assure that ASI will meet your testing needs in a timely manner, **ASI WILL PROCEED WITH THE TESTING** of these samples as directed and comment on the final reports per NELAC requirements. **PLEASE NOTIFY ASI IMMEDIATELY** if you wish to **SUSPEND** analysis, **MAKE ANY CHANGES** to the requested testing services or if the action indicated **IS INCORRECT**.

Sample Integrity Evaluation on Receipt

| | Y | N | N/A | | | Y | N | N/A | |
|----|-------------------------------------|-------------------------------------|--------------------------|---|----|-------------------------------------|-------------------------------------|--------------------------|-------------------------------|
| 1 | <input checked="" type="checkbox"/> | <input checked="" type="checkbox"/> | <input type="checkbox"/> | C-O-C Received w/samples? | 6 | <input checked="" type="checkbox"/> | <input checked="" type="checkbox"/> | <input type="checkbox"/> | Sample Preservation-Temp OK? |
| 2 | <input checked="" type="checkbox"/> | <input checked="" type="checkbox"/> | <input type="checkbox"/> | C-O-C complete with adequate info? | 7 | <input checked="" type="checkbox"/> | <input checked="" type="checkbox"/> | <input type="checkbox"/> | Samples received on ice? |
| 3 | <input checked="" type="checkbox"/> | <input checked="" type="checkbox"/> | <input type="checkbox"/> | C-O-C and samples match (# and descrip.)? | 8 | <input checked="" type="checkbox"/> | <input checked="" type="checkbox"/> | <input type="checkbox"/> | Sample Preservation-pH OK? |
| 4 | <input checked="" type="checkbox"/> | <input checked="" type="checkbox"/> | <input type="checkbox"/> | Custody Seals (if present) intact? | 9 | <input checked="" type="checkbox"/> | <input checked="" type="checkbox"/> | <input type="checkbox"/> | Sample Containers Appropriate |
| 5 | <input checked="" type="checkbox"/> | <input checked="" type="checkbox"/> | <input type="checkbox"/> | Sample Integrity OK? | 10 | <input checked="" type="checkbox"/> | <input checked="" type="checkbox"/> | <input type="checkbox"/> | VOA headspace OK? |
| 11 | <input checked="" type="checkbox"/> | <input checked="" type="checkbox"/> | <input type="checkbox"/> | Dissolved metal samples field filtered and preserved? | | | | | |

Comment: _____

- ☐ ASI Personnel assisted with completion of the C-O-C (in-person or by phone/e-mail).
☐ Additional information supplied w/C-O-C by client.
☐ Samples submitted significantly after (>2 days) sampling, potentially affecting ability to meet hold times.

Comment: _____

Project Management Observations or Discrepancies

- ☐ Insufficient information supplied to determine target analytes required. ASI standard lists will be used.
☐ Special report formats **REQUIRED**. ☐ TRRP ☐ Landfill ☐ Other
☐ Historical project data available for review.
☐ Target analyte list attached.

Comment: _____

Form Sent to Client on: _____ at _____ by ☐ FAX ☐ E-Mail ☐ mail
Client Response Recd.: _____ at _____ by ☐ FAX ☐ E-Mail ☐ VERBAL

Client Response: ☐ Proceed w/analysis ☐ Resample and re-submit

Authorized by (Client Signature): _____ Date _____

[illegible]

APPENDIX B

Removal Action Work Plan Addendum No. 3

Falcon Refinery Superfund Site Ingleside San Patricio County, Texas TXD 086 278 058

Prepared for

**National Oil Recovery Corporation
3717 Bowne Street
Flushing, NY 11354**

Prepared by:



**TRC Environmental Corporation
505 East Huntland Drive, Suite 250
Austin, Texas 78752**

May 2011

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FIGURES

Figure 1 Area Map

Figure 2 Above Ground Storage Tank Map

1. INTRODUCTION

This report describes past activities performed by National Oil Recovery Corporation (NORCO) as well as upcoming activities associated with the EPA approved Removal Action Work Plan, which was directed by the *Administrative Order on Consent (AOC)*, *CERCLA Docket No 06-04-04* at the Falcon Refinery in Ingleside, Texas.

Provided below are excerpts from the AOC that are pertinent to the scope of work. Each of the required items in the section termed “The Work” will be described in terms of work already completed and work yet to be completed.

Actions associated with the Removal Action were performed under the EPA approved Removal Action Work Plan, dated July 20, 2004.

WORK TO BE PERFORMED

The Respondent shall perform, at a minimum, all actions necessary to implement the Order of Work. The actions to be implemented generally include, but are not limited to, the following:

The Work

The intent of this action is to remove and dispose of source materials from tanks and other miscellaneous containers, equipment, piping and buildings. This also includes the removal of any source materials in piping associated with transfer or materials to the facility docks or former facility docks. As a result, it may be necessary to demolish or otherwise remove some tanks, piping, equipment, and buildings in order to effectuate this activity. This action may also include the consolidation of materials in onsite tankage for future disposal. If the Respondent elects to consolidate and temporarily store these materials, Respondent must comply with all applicable laws including storage and spill prevention regulations.

The work includes:

- *Asbestos Inspection and Abatement: The Respondent shall comply with applicable law(s) to address any asbestos and coordinate its handling appropriately with any demolition activities. Therefore, it will be necessary for the Respondent to conduct an asbestos inspection and make appropriate notifications for the conduct of such demolition and asbestos abatement activities as required by applicable law(s).*
- *Assessment and Removal of Hazardous Substances, or Pollutants or Contaminants: The Respondent shall conduct tests and properly classify the wastes for appropriate disposal or recycling.*
- *Decontamination of Containers, Equipment, Piping, and Buildings: The Respondent shall decontaminate all containers, equipment, piping, and buildings to the extent necessary for the purpose of recycling, reuse, or disposal.*

- *Removal of Containers, Equipment, Piping, and other Contaminated Items: The Respondent shall recycle or dispose of containers, equipment, piping, and other potentially contaminated items in accordance with applicable law(s). The metal debris associated with the removal of the containers, equipment, piping, and other items should be recycled to the extent practical.*
- *Consolidation, Removal and Disposal of Grossly Contaminated Soil: The Respondent shall consolidate and then treat or dispose of visibly contaminated surface soils identified during the conduct or resulting from the conduct of this action.*

2. COMPLETED ACTIVITIES

Described in this section are the completed activities associated with “The Work”.

2.1. Asbestos Inspection and Abatement

Asbestos sampling, which was performed during August 2004 and reported in the September 2004 Monthly Progress Report, indicated that minimal asbestos was present at the site. Detected asbestos containing materials (ACM) were limited to gaskets associated with various pipeline connections. No friable asbestos materials were present.

2.2. Assessment and Removal of Hazardous Substances, or Pollutants or Contaminants

During September 2004 the contents of all the above ground storage tanks (AST) were gauged to determine the volume of waste and sampled to characterize the waste. An estimated 6.8 million gallons of liquid waste that required disposal was measured. In addition to the liquid waste 62,000 gallons of sludge was measured.

Analytical sampling indicated varied waste streams and compatibility testing was performed of the waste to ensure safe disposal.

Based on the characteristics of the waste deep well injection at the Texas Molecular facility in Corpus Christi was selected. Three tanker trucks a day made three trips each from the refinery to the Texas Molecular facility carrying liquid waste, which resulted in the disposal of 7,774,721 gallons of waste. The waste disposal volume was higher than the estimated amount due to rainfall entering the tanks during disposal operations.

2.3. Decontamination and Removal of Containers, Equipment, Piping, and Buildings

From October 2004 through February 2005 the onsite buildings, abandoned drums and equipment were described, decontaminated, characterized and properly disposed. Results included the recycling of 67,840 pounds of metal, 10 fire extinguishers, and 403 gallons of oil and filters.

Items that couldn't be recycled were disposed, which resulted in the disposal of 16,651 gallons of waste oil.

When oil was discovered leaking in the wetlands adjacent to the refinery, ten pipelines were excavated, cut at five locations and jetted clean prior to having steel caps welded on the pipelines.

2.1.Consolidation, Removal and Disposal of Grossly Contaminated Soil

During September 2004 measurements of grossly contaminated soil based on the surface extent and depth of visibly impacted soil was estimated to be approximately 6,000 cubic yards. Based on the amount of grossly contaminated soil NORCO proposed, to the EPA, the possible construction of bioremediation cells to treat the soil on site rather than disposing of the soil at a hazardous waste facility.

During December 2004 and February 2005, 55 cubic yards of grossly contaminated soil was disposed at the US Ecology Texas Facility in Robstown, with EPA's approval.

The remainder of the grossly contaminated soil cannot be excavated until the sludge is removed from the above ground storage tanks.

3. UPCOMING ACTIVITIES

Described in this section are the upcoming activities associated with “The Work”.

3.1. Asbestos Inspection and Abatement

Based on the results of the asbestos sampling performed during 2004, which indicated no friable asbestos, no asbestos inspection or abatement is anticipated. If any asbestos containing materials (ACM) are detected or if a material appears to be ACM, testing will be performed and appropriate measures will be taken.

3.2. Assessment and Removal of Hazardous Substances, or Pollutants or Contaminants

As of September 2009 all liquid waste was removed from all of the above ground storage tanks. However, sludge was measured and remained in Tanks 7, 10, 26, 27 and 30.

Upon approval of the commencement of work by NORCO the initial action will be the measurement of the contents of all tanks and vessels. Tanks leased to Superior Crude Gathering (Superior), which includes Tanks 13, 15 and 16 will not be inspected as they are either in use or are in the process of being repaired. An additional unnamed and unnumbered tank, noted by the EPA as leaking during an inspection will also be evaluated.

After the measurement of the sludge NORCO anticipates mobilizing a centrifuge to separate liquid waste from solid waste. The last estimate of the volume of sludge indicated that there would be approximately 180 tons of solid waste and 15,000 gallons of liquid waste that will need disposal.

When the materials are separated they will be stored in appropriate containers (frac tanks or lined roll off boxes) pending characterization and disposal. The EPA will be notified of any disposal plans and no disposal will take place prior to EPA approval.

3.3. Decontamination and Removal of Containers, Equipment, Piping, and Buildings

When the contents of the tanks are removed a determination will be made about the usability of each tank. If a tank is to be used in the future the inside of the tank will be either steam cleaned or sand blasted and a full American Petroleum Institute (API) 653 internal/external (out-of-service) inspection will be conducted by a properly certified API 653 tank inspector.

For tanks that remain in service repairs will begin immediately to prevent the collection of rainwater in tanks due to leaking roofs.

If a tank is to be razed then the tank will still be cleaned to the level necessary for disposal. The EPA will be notified of the future of each tank and will be allowed to observe the cleaning of the tanks.

Prior to any tank parts leaving the site the EPA will be notified.

Currently there are no plans to remove any of the former refining equipment or to construct any new refining equipment. Prior to any construction the EPA will be notified. Any removed metal will be recycled.

3.4. Consolidation, Removal and Disposal of Grossly Contaminated Soil

During a site inspection in March 2011 there was no grossly contaminated soil observed at the main portion of the refinery or around the storage tanks. The inspection did not include the north property as a result there may be grossly contaminated soil on that property.

On February 9, 2010 Superior, which leases several of the above ground tanks at the site, had a release from Tank 13. Approximately 22,000 barrels of crude oil was released from the tank and onto the site. Remediation efforts, using pumps and vacuum trucks were successful in the recovery of much of the crude oil.

Superior used heavy equipment to scrape off the impacted soil (grossly contaminated), which is still stockpiled at the site pending regulatory approval for disposal. The results were described in the *Site Investigation Report and Remediation Plan, Superior Crude Gathering, Inc, Crude Oil Spill, Ingleside, Texas* dated July 23, 2010 by Pastor, Behling & Wheeler.

Upon approval of the commencement of work by NORCO an assessment will be made of the volume of grossly contaminated soil and the EPA will be notified. Based on the volume of soil NORCO will either propose the construction of a bioremediation cell to treat the soil or will propose disposal at an appropriate facility. No soil will be moved without EPA approval.

After each tank is cleaned as noted in section 3.3 a determination will be made about the re-use of the tank. If the tank is removed then the soil underneath the tank will be checked for visual contamination. If the tank is not taken out of service then the bottom of the tank will be thoroughly inspected to insure that there are no cracks, holes or areas of weakness that could have resulted in the release of contamination below the tank. If inspection reveals damage to the tank bottom then the tank bottom will be removed and any visually contaminated soil will be removed prior to the repair or replacement of the tank bottom.

Any excavated grossly contaminated soil will be replaced by clean fill brought to the site.

4. SCHEDULE

NORCO will begin work when notified by the EPA that actions may commence. The initial tank inspection is expected to take one week to determine the volume of waste in the tanks. Once the volume of waste is determined equipment likely including a centrifuge, frac tanks, pumps, vacuum trucks and associated equipment will be mobilized to the site to begin the removal of the contents of the tanks that contain waste.

As noted previously when the tanks are emptied the tanks will be cleaned to the appropriate level and any grossly contaminated soil around the tanks or around the site will be excavated and either treated on site or disposed of properly.

All work associated with the Removal Action will be completed by December 31, 2011 and all tankage will be cleaned and gas freed no later than August 1, 2011.